

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
 Data File : BF092571.D
 Acq On : 27 Jan 2017 12:27
 Operator : SJ/MA
 Sample : I1353-07
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-02

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-BF012417.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	3.195	94	97	99	rBV	45113	83959	1.47%	0.086%
2	5.161	266	269	271	rBV	536319	708478	12.41%	0.728%
3	5.219	272	274	278	rVB2	68564	133490	2.34%	0.137%
4	5.447	290	294	296	rBV	51765	77908	1.36%	0.080%
5	5.481	296	297	300	rVB	67626	77323	1.35%	0.079%
6	5.584	304	306	308	rVV	2687603	2433229	42.62%	2.499%
7	5.687	313	315	317	rBV	255021	278866	4.88%	0.286%
8	5.744	318	320	324	rVB	183292	242507	4.25%	0.249%
9	5.847	327	329	330	rBV2	111876	109406	1.92%	0.112%
10	5.881	330	332	334	rVB	101526	119899	2.10%	0.123%
11	6.007	338	343	345	rBV4	55522	114972	2.01%	0.118%
12	6.133	348	354	355	rBV3	137600	211090	3.70%	0.217%
13	6.167	355	357	358	rBV	58777	72940	1.28%	0.075%
14	6.201	358	360	361	rBV	153616	129902	2.28%	0.133%
15	6.224	361	362	365	rVB2	110036	175067	3.07%	0.180%
16	6.282	365	367	371	rVB4	68273	136226	2.39%	0.140%
17	6.384	374	376	378	rBV2	62243	70839	1.24%	0.073%
18	6.419	378	379	382	rVB	197471	219222	3.84%	0.225%
19	6.464	382	383	384	rBV	96959	75220	1.32%	0.077%
20	6.544	384	390	391	rBV3	62597	153272	2.68%	0.157%
21	6.567	391	392	394	rBV	261899	228180	4.00%	0.234%
22	6.624	394	397	400	rBV	1163735	1782619	31.22%	1.831%
23	6.670	400	401	403	rVB	96706	101842	1.78%	0.105%
24	6.750	405	408	410	rBV	3478606	4190443	73.39%	4.304%
25	6.819	411	414	415	rBV	263701	332942	5.83%	0.342%
26	6.842	415	416	419	rVB2	120873	163743	2.87%	0.168%
27	6.922	419	423	425	rBV4	416100	677492	11.87%	0.696%
28	6.967	425	427	429	rVB2	1038004	1094015	19.16%	1.124%
29	7.036	432	433	438	rVB2	199083	281411	4.93%	0.289%
30	7.127	438	441	446	rBV2	3508505	5155377	90.29%	5.295%
31	7.219	446	449	451	rBV2	787209	778423	13.63%	0.799%
32	7.253	451	452	454	rVB2	105931	125806	2.20%	0.129%
33	7.310	454	457	461	rVB2	620052	1456917	25.52%	1.496%
34	7.379	461	463	464	rBV2	280124	362841	6.35%	0.373%

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 Sampling : 1
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 Stop Thrs : 0

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012417.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	7.413	464	466	470	rBV3	422069	804685	14.09%	0.826%
36	7.470	470	471	472	rBV	261140	266305	4.66%	0.274%
37	7.516	472	475	476	rBV2	1783499	2835915	49.67%	2.913%
38	7.550	476	478	480	rVB	2901850	3373648	59.09%	3.465%
39	7.596	480	482	483	rVB2	266395	249893	4.38%	0.257%
40	7.619	483	484	487	rVB3	698961	798975	13.99%	0.821%
41	7.665	487	488	489	rBV	463569	407160	7.13%	0.418%
42	7.745	493	495	497	rVB	1300924	1405992	24.62%	1.444%
43	7.802	497	500	501	rBV	592977	747440	13.09%	0.768%
44	7.847	501	504	508	rVB3	992792	1289590	22.59%	1.324%
45	7.939	508	512	513	rBV2	808295	2096638	36.72%	2.153%
46	7.973	513	515	516	rVB	386680	309452	5.42%	0.318%
47	7.996	516	517	519	rBV2	1019721	1434326	25.12%	1.473%
48	8.030	519	520	522	rVB2	1055144	1306303	22.88%	1.342%
49	8.099	522	526	528	rBV4	701992	2078223	36.40%	2.134%
50	8.133	528	529	530	rVB	549874	376939	6.60%	0.387%
51	8.213	534	536	537	rVB2	552942	547024	9.58%	0.562%
52	8.270	537	541	544	rBV3	2350633	5709645	100.00%	5.864%
53	8.385	548	551	553	rBV3	418023	768320	13.46%	0.789%
54	8.465	555	558	561	rBV2	1370882	2808736	49.19%	2.885%
55	8.522	561	563	566	rBV4	590364	816540	14.30%	0.839%
56	8.613	569	571	572	rVB2	776702	865976	15.17%	0.889%
57	8.636	572	573	576	rVB3	864125	1349269	23.63%	1.386%
58	8.727	576	581	582	rBV3	1218271	2788254	48.83%	2.864%
59	8.773	583	585	588	rVB4	803096	1247225	21.84%	1.281%
60	8.819	588	589	591	rVB2	1031151	820428	14.37%	0.843%
61	8.865	591	593	595	rVB2	832850	1275890	22.35%	1.310%
62	8.933	595	599	601	rBV2	1739366	3263563	57.16%	3.352%
63	8.967	601	602	604	rBV2	691374	720941	12.63%	0.740%
64	9.025	605	607	608	rVB	520208	403435	7.07%	0.414%
65	9.059	608	610	611	rBV2	701950	750725	13.15%	0.771%
66	9.093	611	613	616	rBV3	964177	1882074	32.96%	1.933%
67	9.162	616	619	621	rBV4	744465	1920497	33.64%	1.972%
68	9.299	628	631	634	rBV4	848770	2415752	42.31%	2.481%
69	9.356	634	636	640	rVB	3980020	5230297	91.60%	5.372%
70	9.413	640	641	643	rBV	1725861	2506396	43.90%	2.574%
71	9.539	649	652	653	rBV3	916341	1584825	27.76%	1.628%

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72	9.825	671	677	679	rBV5	1090182	3764460	65.93%	3.866%
73	9.893	681	683	685	rVB	1397887	1364573	23.90%	1.401%
74	9.973	688	690	693	rBV3	874683	1553408	27.21%	1.595%
75	10.042	693	696	698	rVV2	1300066	2189619	38.35%	2.249%
76	10.099	699	701	702	rVB	837047	952711	16.69%	0.978%
77	10.385	723	726	728	rBV	806469	1951399	34.18%	2.004%
78	13.128	964	966	969	rVB	1643211	2026612	35.49%	2.081%
79	14.179	1056	1058	1061	rVB	645866	836342	14.65%	0.859%
80	15.677	1186	1189	1192	rVB	658434	913758	16.00%	0.938%

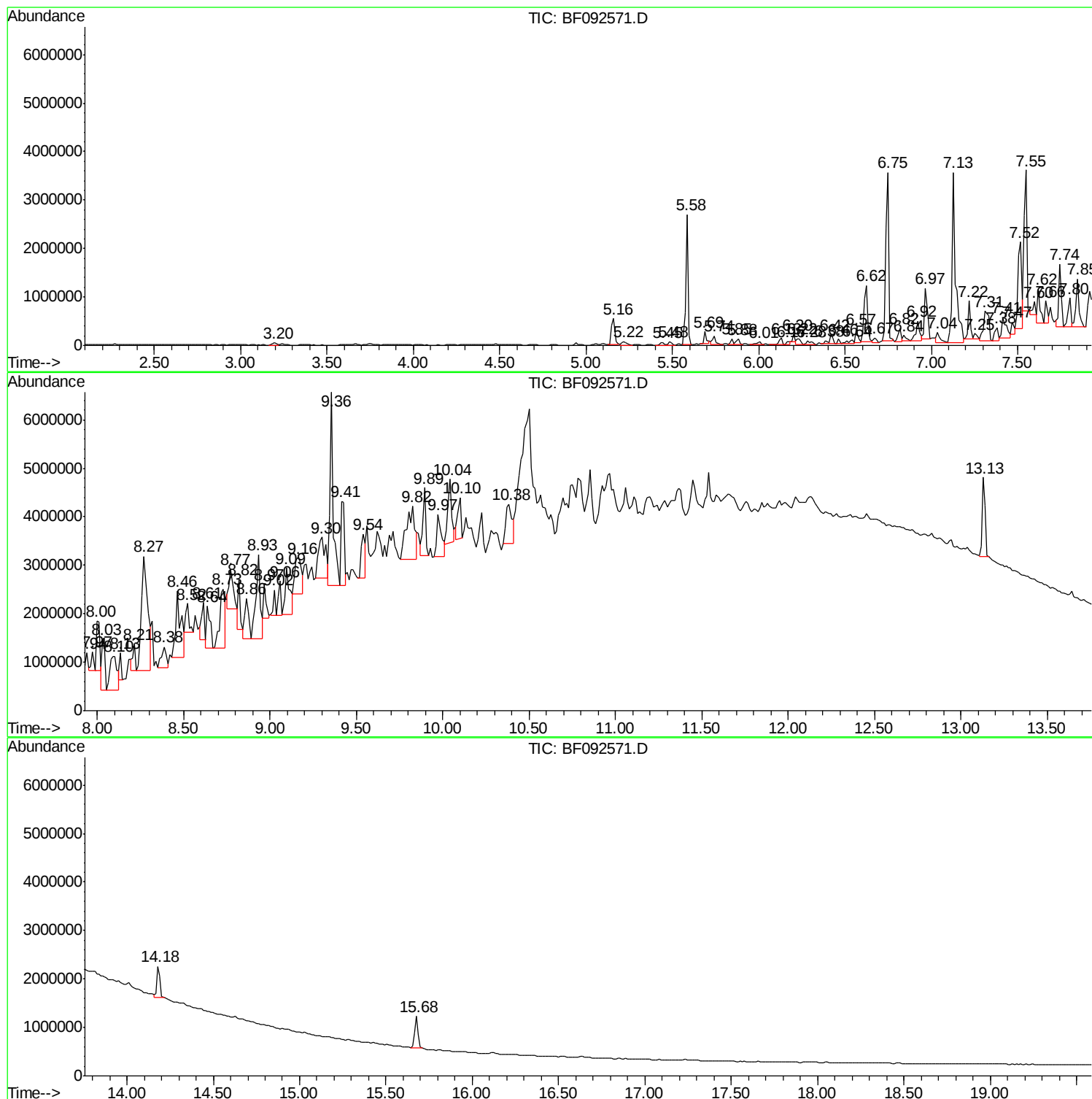
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Instrument :
BNA_F
ClientSampled :
W-02

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
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Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-02

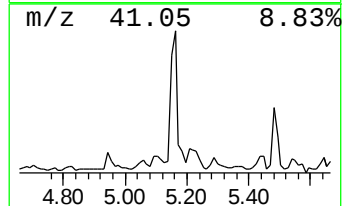
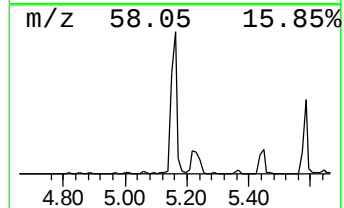
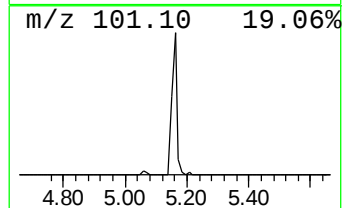
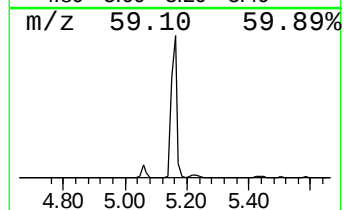
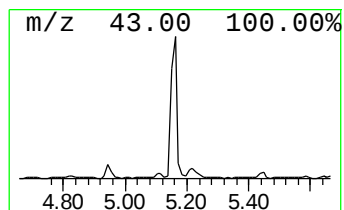
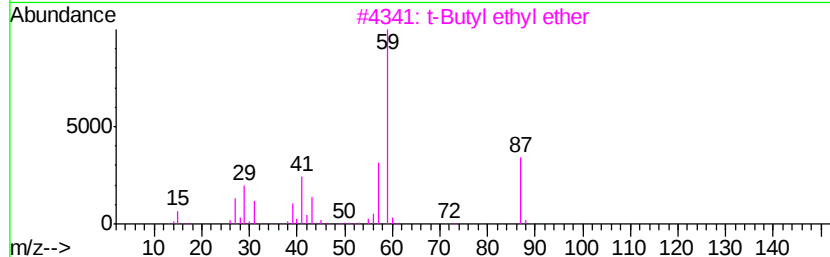
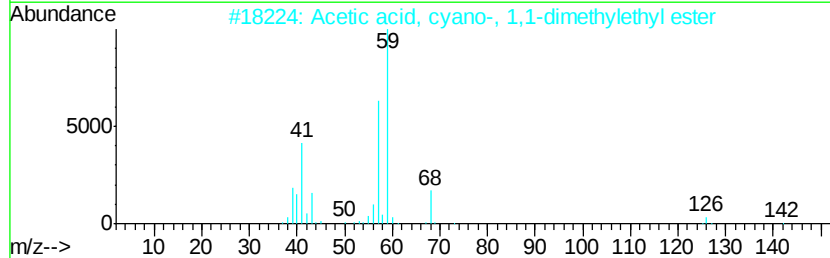
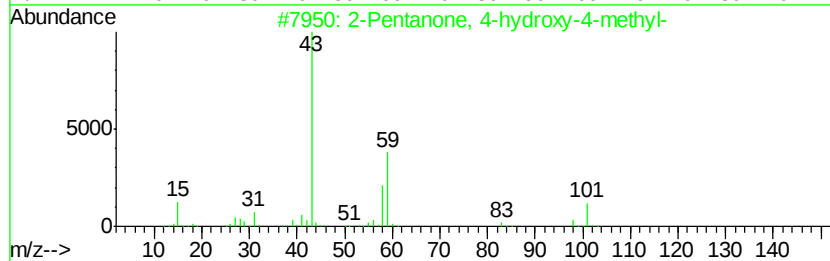
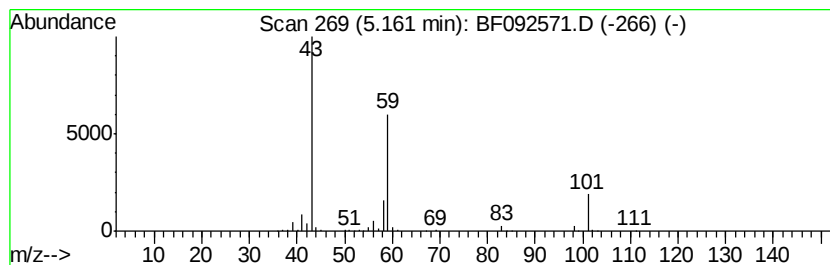
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	12.95 ng	708478	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		t-Butyl ethyl ether	102	C6H14O	1000118-18-4	9
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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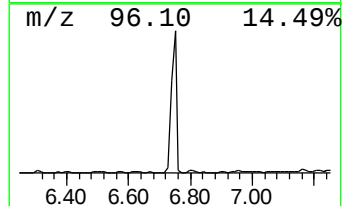
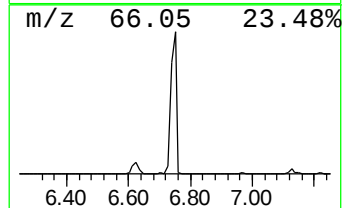
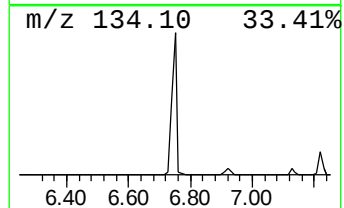
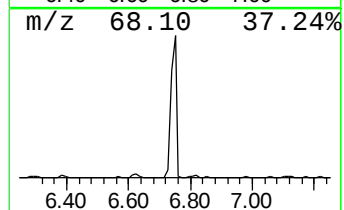
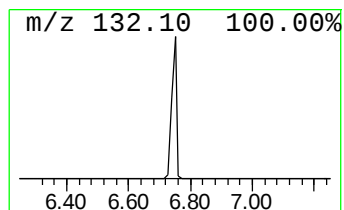
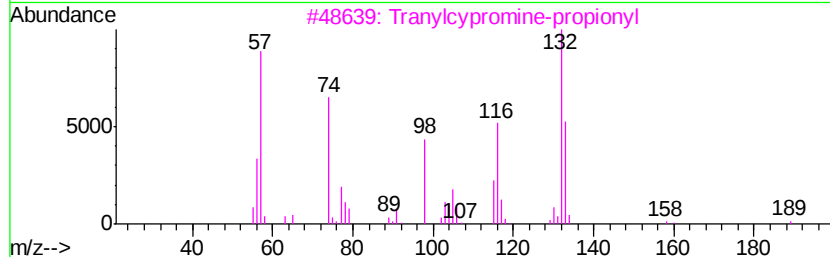
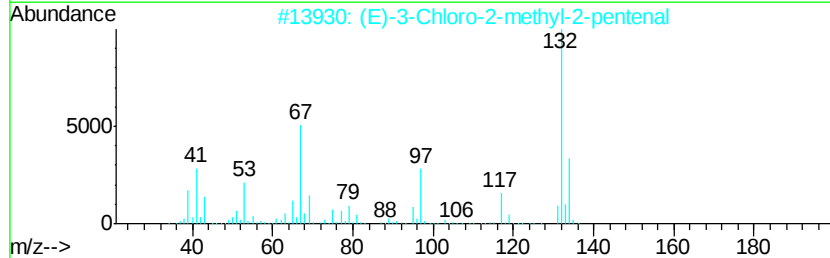
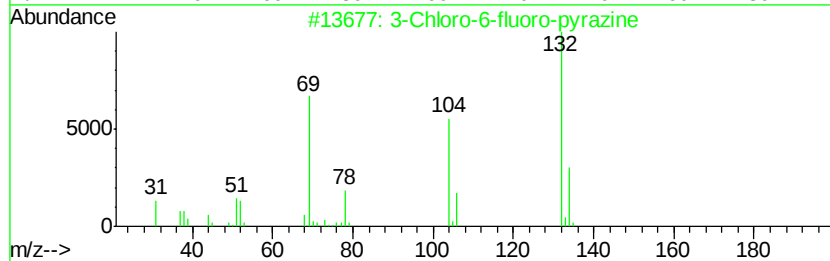
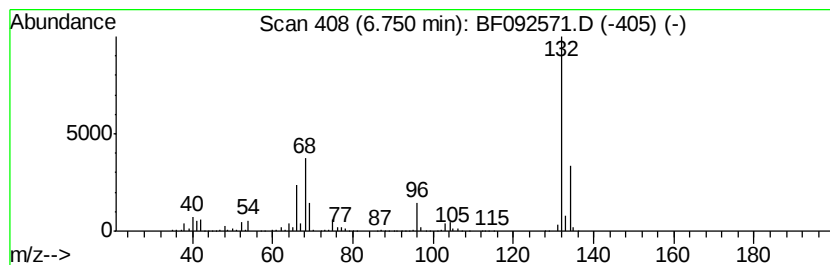
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.75 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	76.61 ng	4190440	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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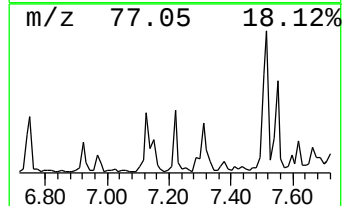
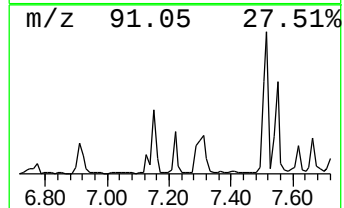
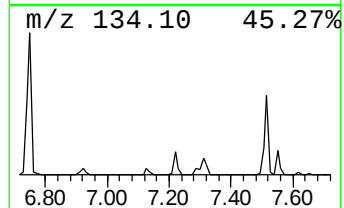
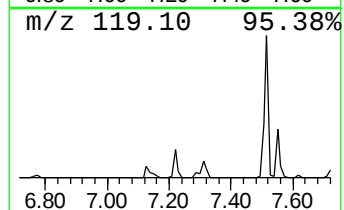
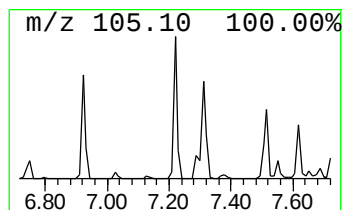
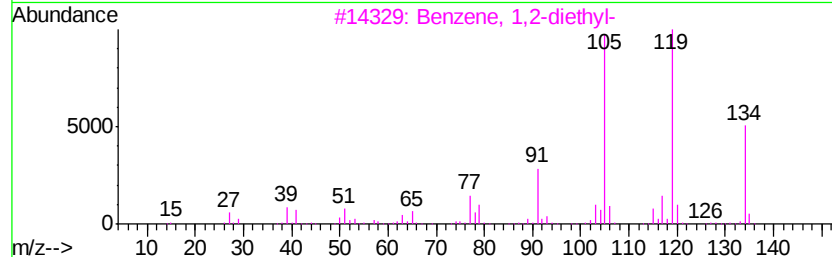
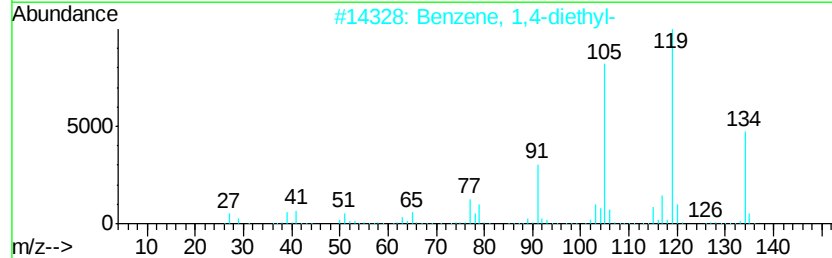
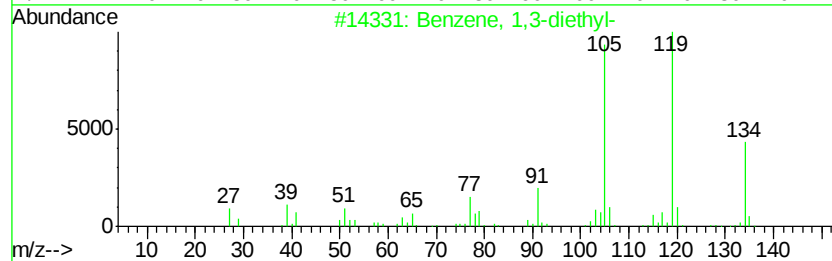
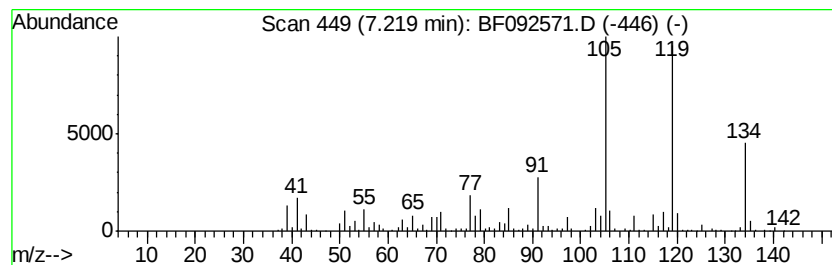
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	14.23 ng	778423	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87



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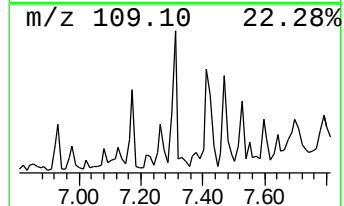
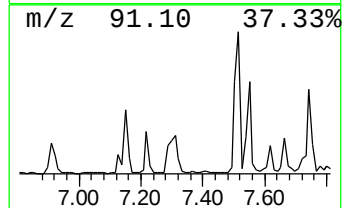
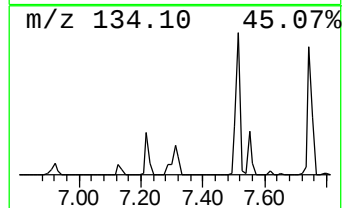
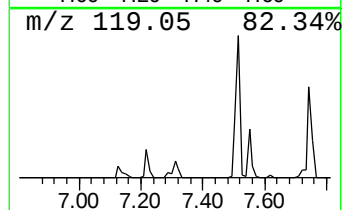
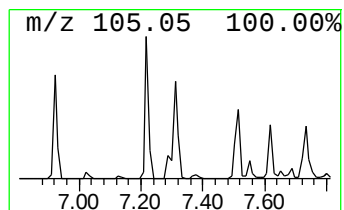
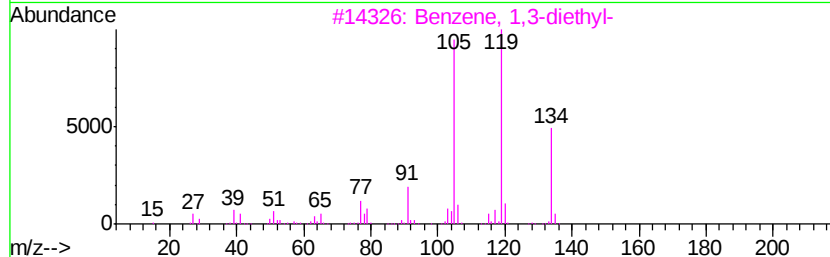
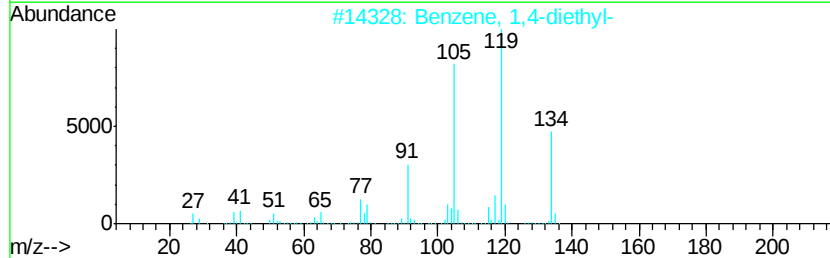
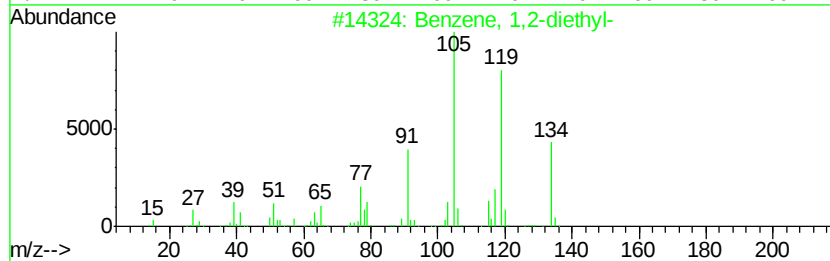
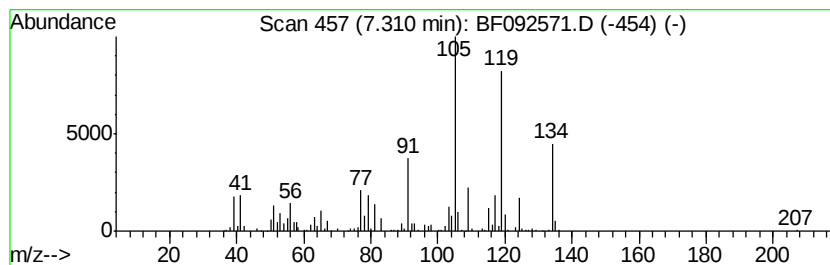
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,2-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.31	26.63 ng	1456920	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	83
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	70
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	68



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-02

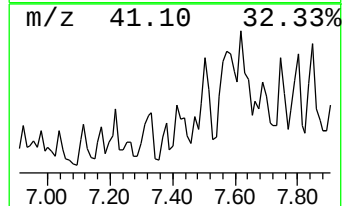
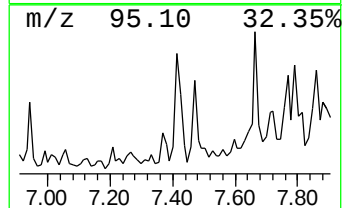
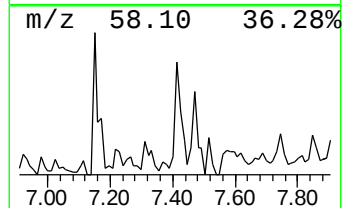
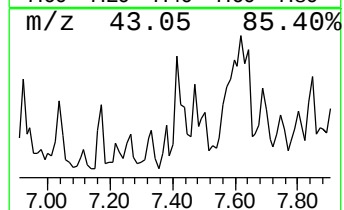
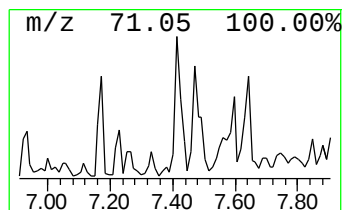
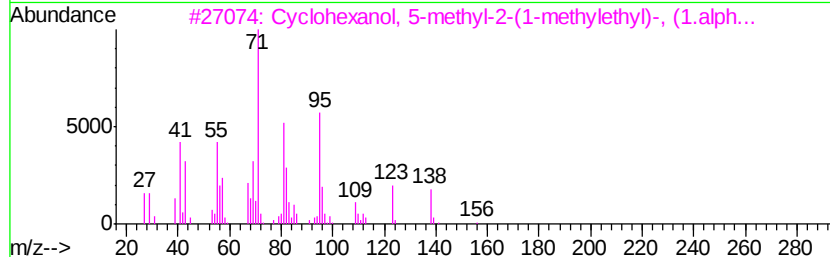
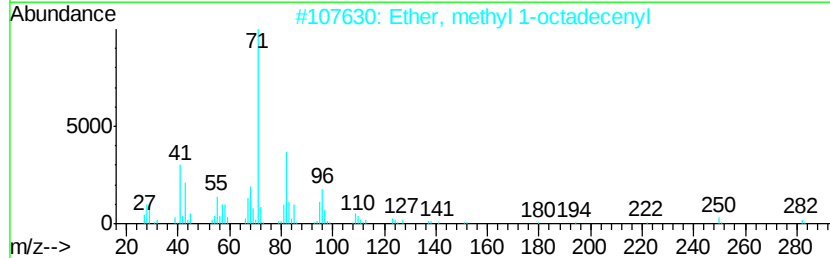
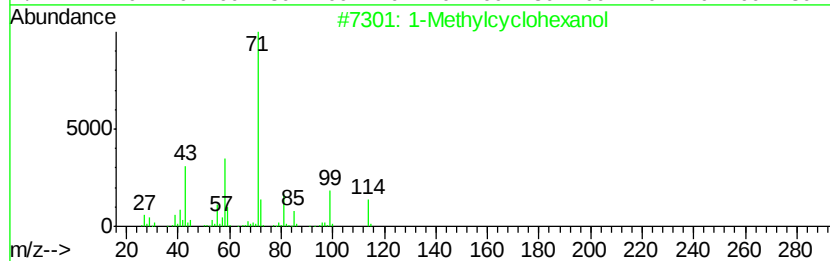
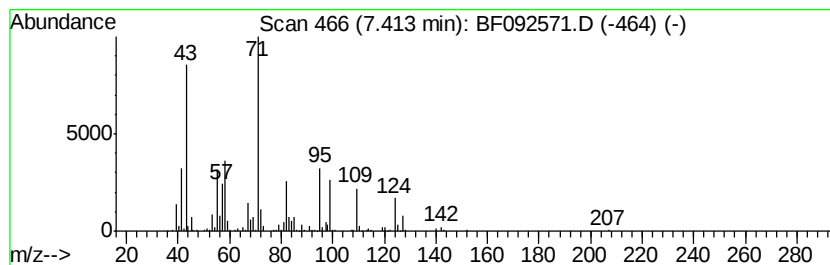
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown7.41 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.41	14.71 ng	804685	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methylcyclohexanol	114	C7H14O	000590-67-0	43
2		Ether, methyl 1-octadecenyl	282	C19H38O	026537-06-4	38
3		Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	000491-02-1	38
4		Cyclooctane, methoxy-	142	C9H18O	013213-32-6	38
5		Oxirane, butyl-	100	C6H12O	001436-34-6	37



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
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Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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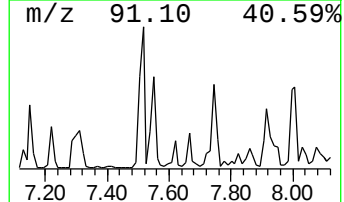
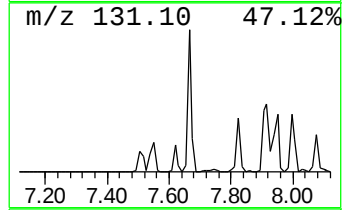
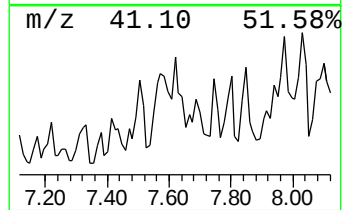
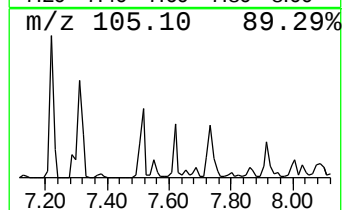
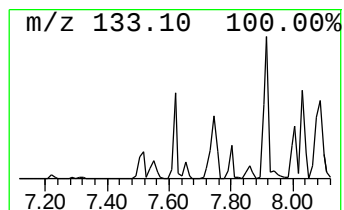
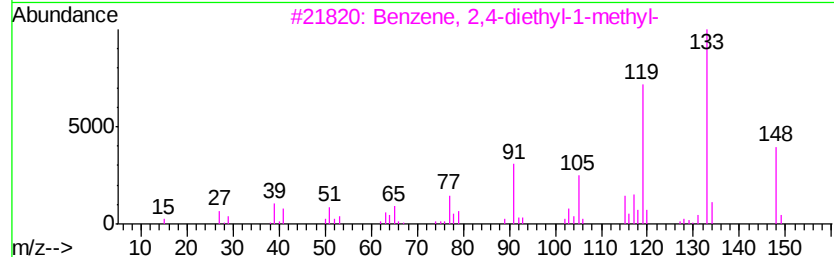
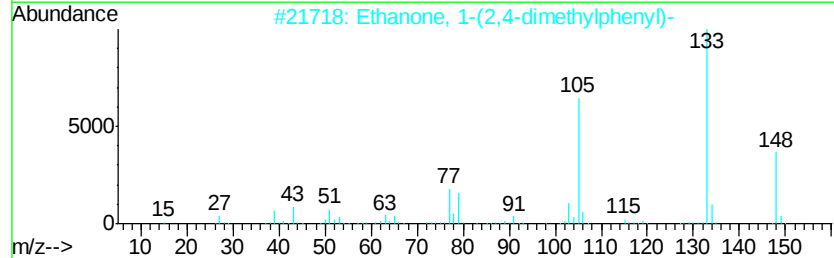
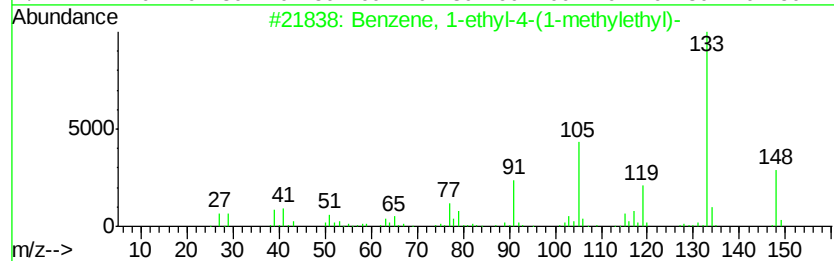
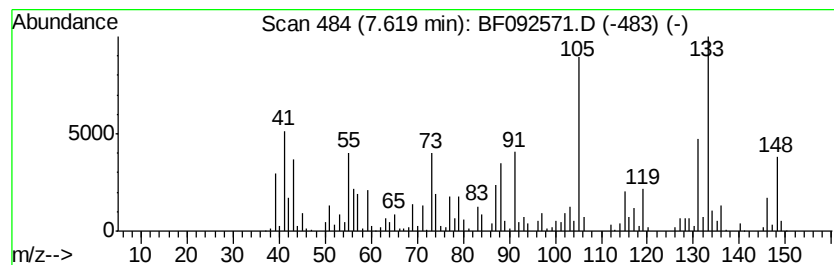
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-4-(1-methylethyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	14.61 ng	798975	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	50
2		Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	46
3		Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	46
4		Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	46
5		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	46



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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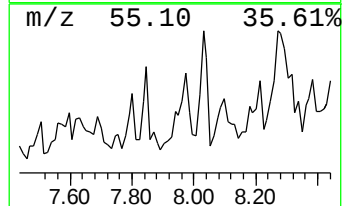
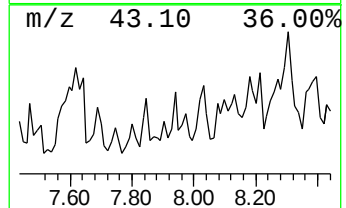
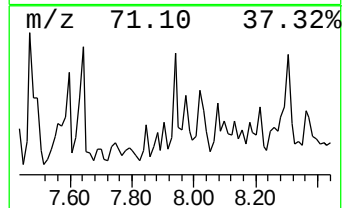
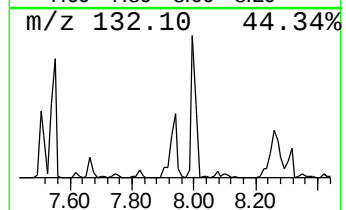
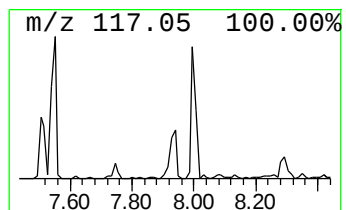
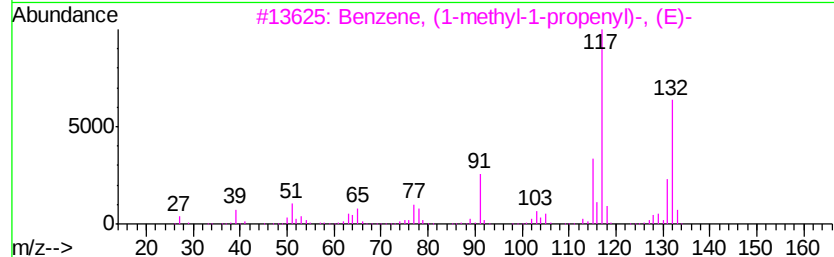
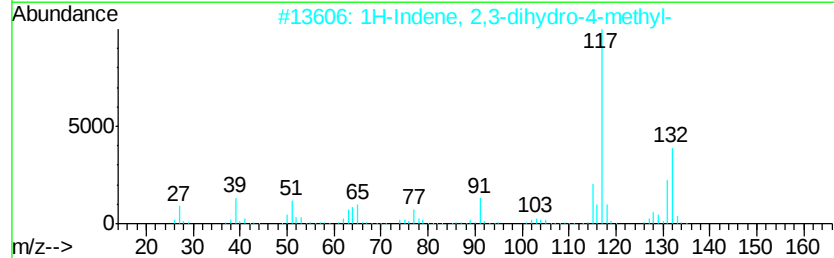
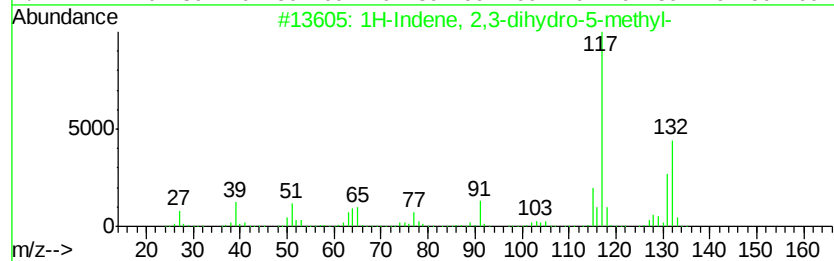
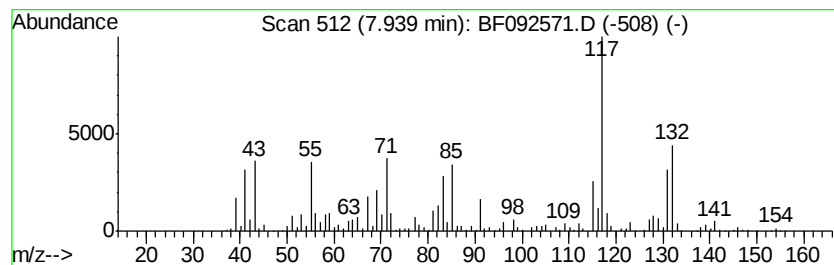
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.94	7.34 ng	2096640	Naphthalene-d8	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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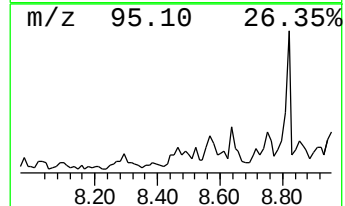
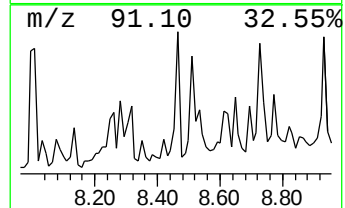
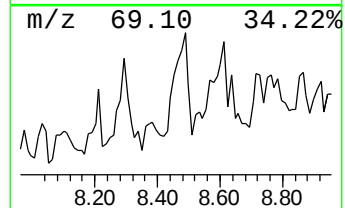
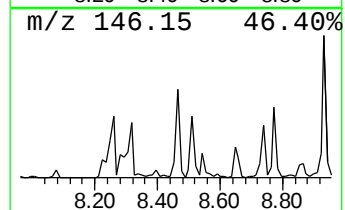
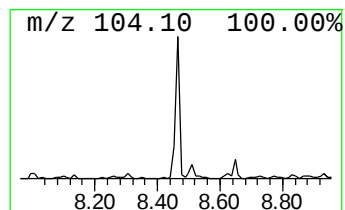
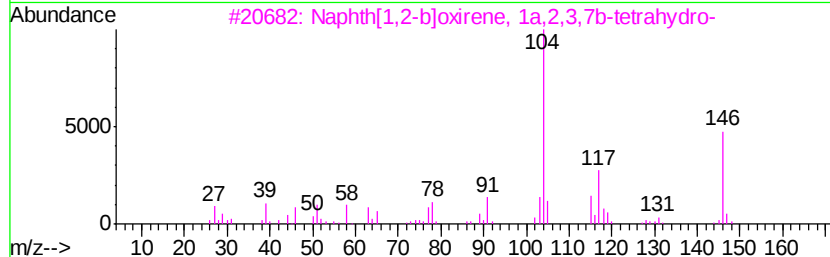
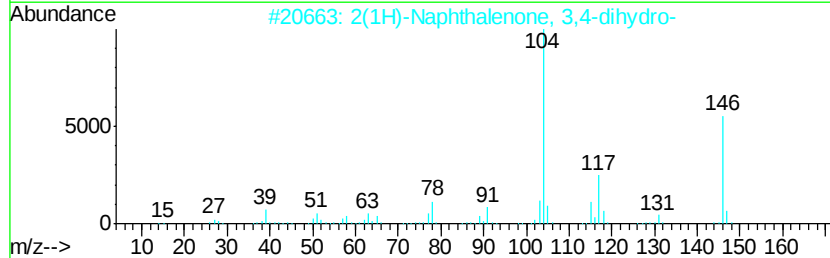
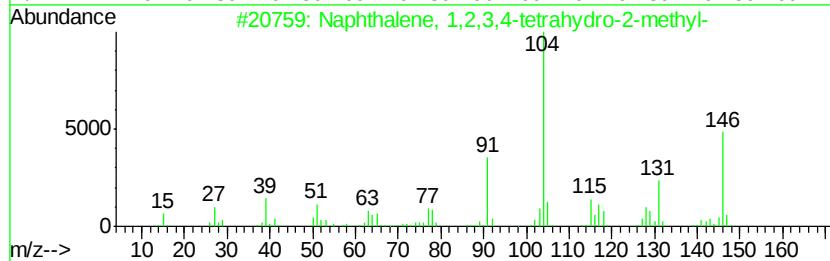
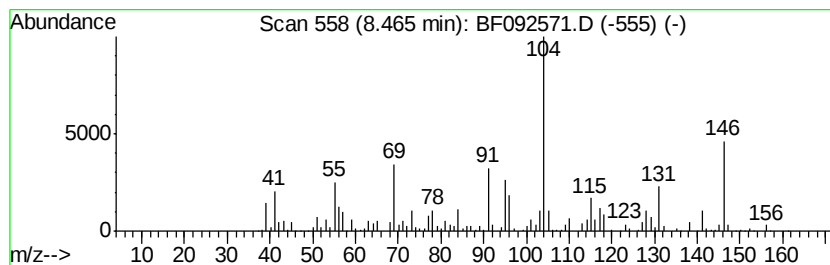
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.46	9.84 ng	2808740	Naphthalene-d8	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	89
2		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	49
3		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	49
4		2,6-Difluorobenzoic acid, 2-phen...	262	C15H12F2O2	1000293-48-2	35
5		Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-02

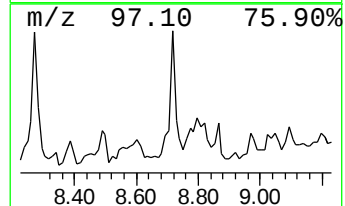
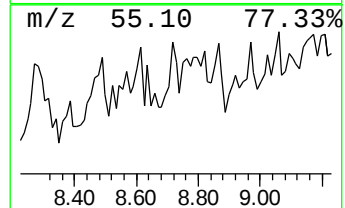
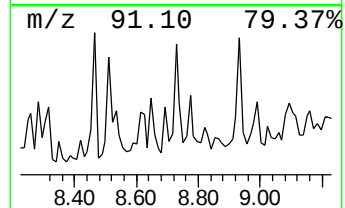
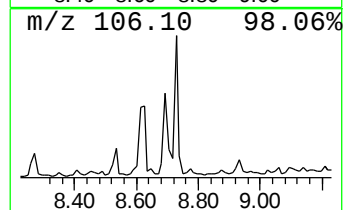
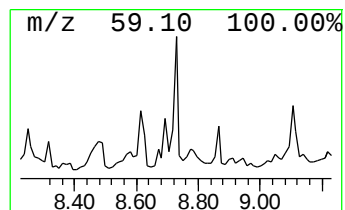
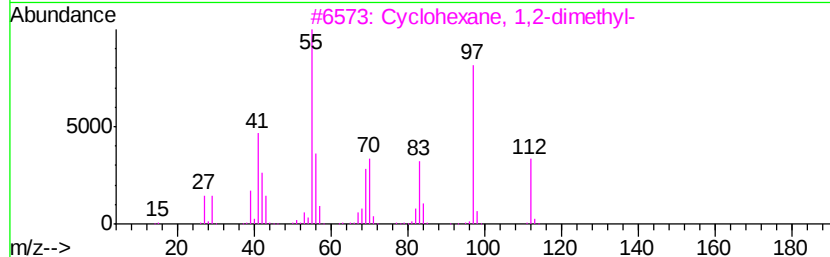
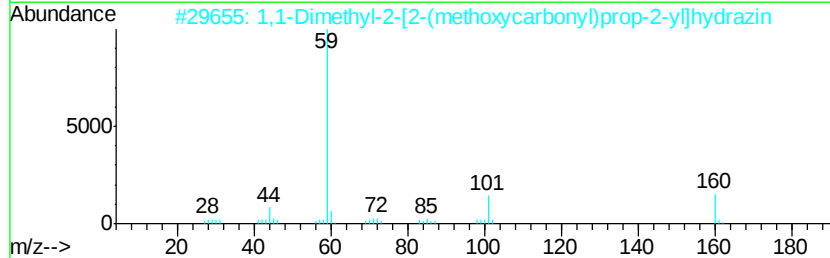
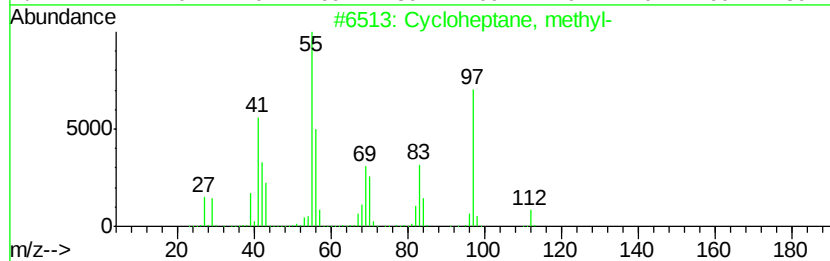
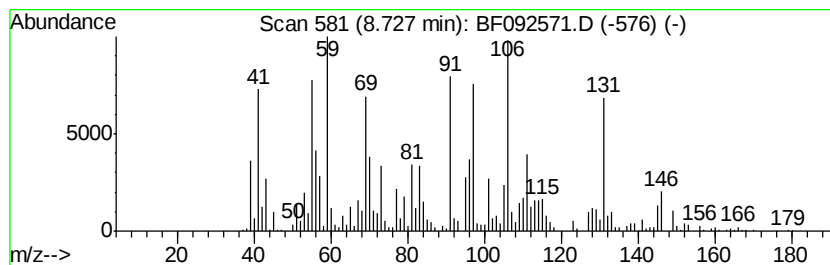
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown8.73 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.73	9.77 ng	2788250	Napthalene-d8	8.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloheptane, methyl-	112	C8H16	004126-78-7	25
2			1,1-Dimethyl-2-[2-(methoxycarbon...	160	C7H16N2O2	097812-30-1	15
3			Cyclohexane, 1,2-dimethyl-	112	C8H16	000583-57-3	15
4			Propanenitrile, 3-(phenylamino)-	146	C9H10N2	001075-76-9	11
5			Ethanol, 2-(phenylamino)-	137	C8H11NO	000122-98-5	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
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Acq On : 27 Jan 2017 12:27
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Misc :
ALS Vial : 56 Sample Multiplier: 1

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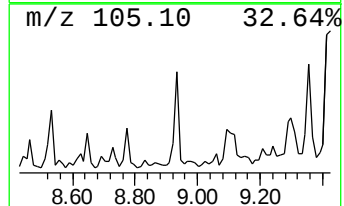
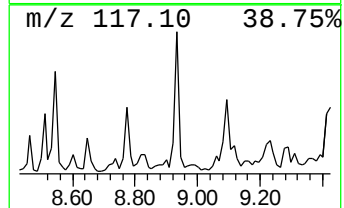
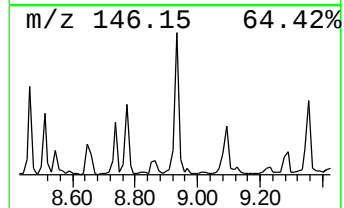
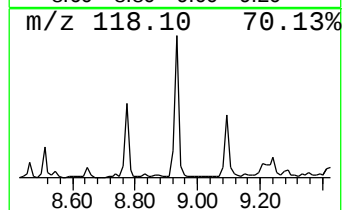
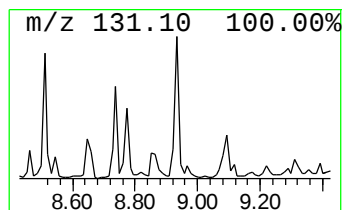
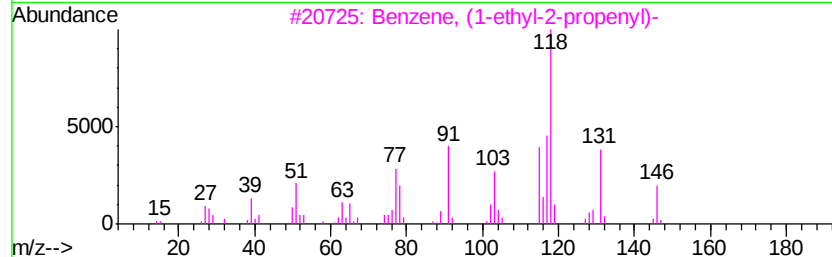
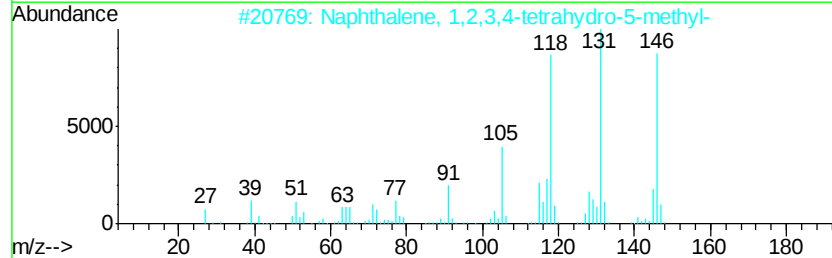
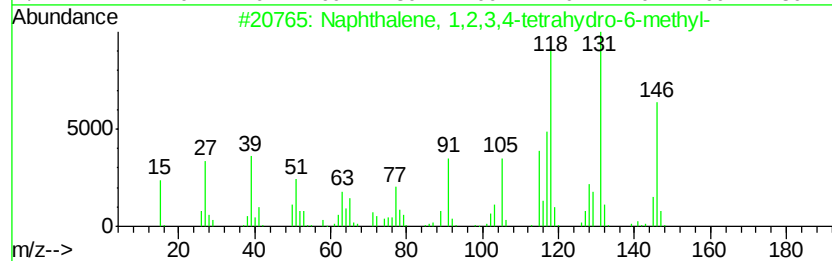
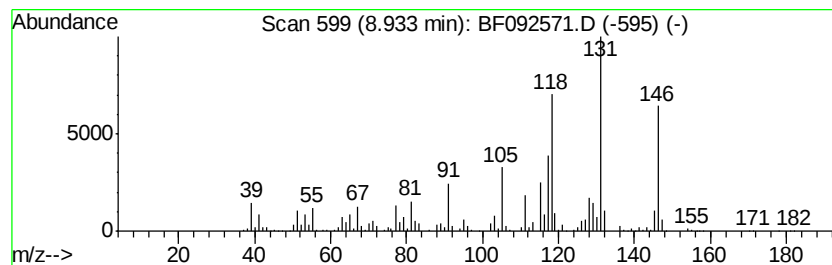
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,2,3,4-tetra... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.93	11.43 ng	3263560	Naphthalene-d8	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3		Benzene, (1-ethyl-2-propenyl)-	146	C11H14	019947-22-9	81
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	70
5		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
W-02

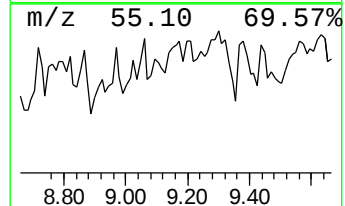
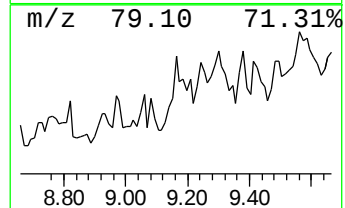
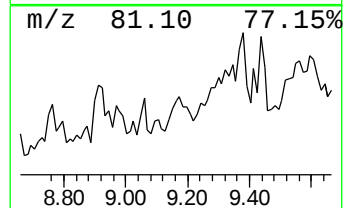
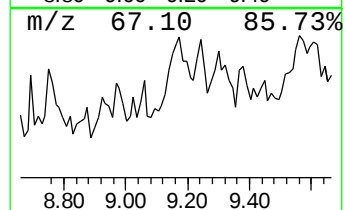
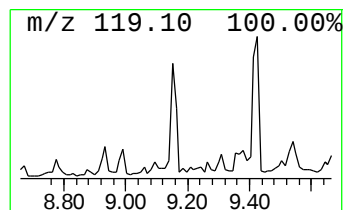
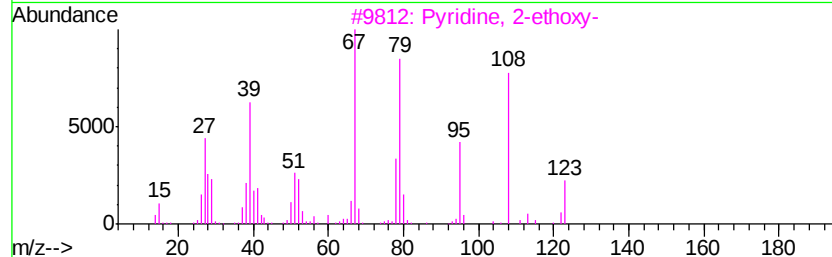
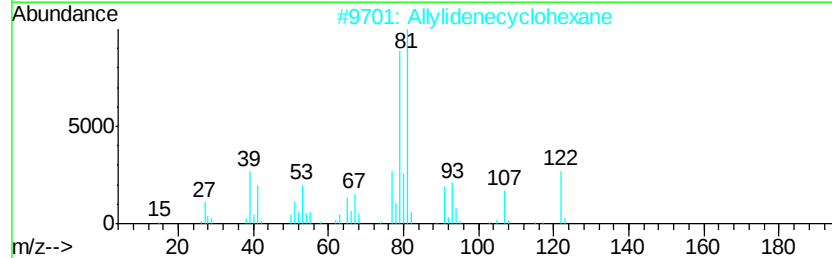
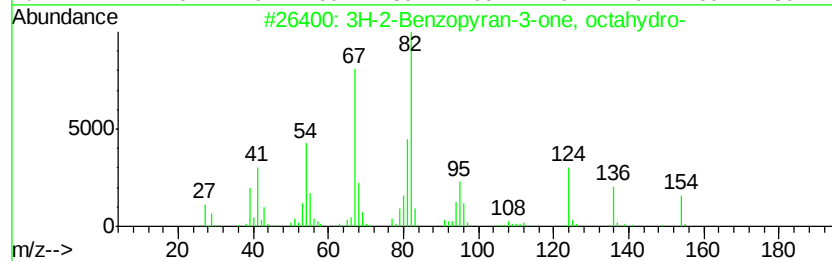
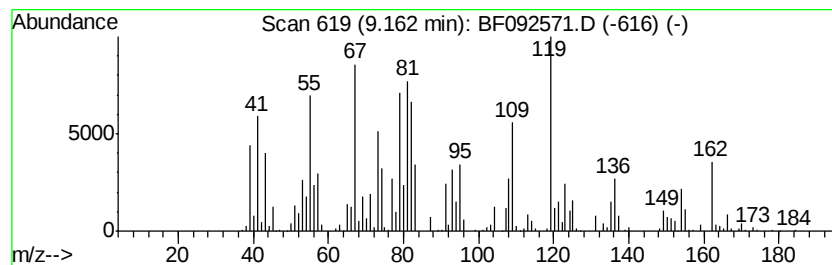
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown9.16 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.16	17.54 ng	1920500	Acenaphthene-d10	10.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3H-2-Benzopyran-3-one, octahydro-	154	C9H14O2	021962-62-9	25
2		Allylidene cyclohexane	122	C9H14	005664-10-8	25
3		Pyridine, 2-ethoxy-	123	C7H9NO	014529-53-4	18
4		1,6-Cyclodecadiene	136	C10H16	007049-13-0	18
5		Cis-3-methyl-exo-tricyclo[5.2.1....	150	C11H18	1000215-28-9	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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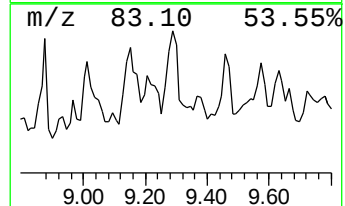
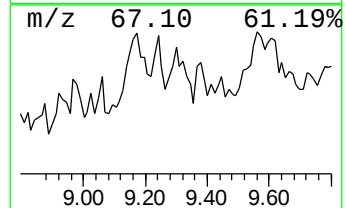
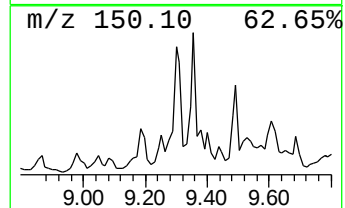
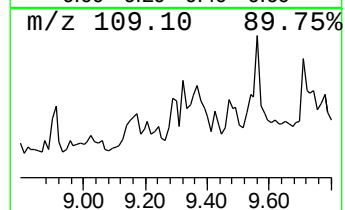
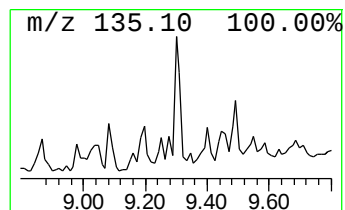
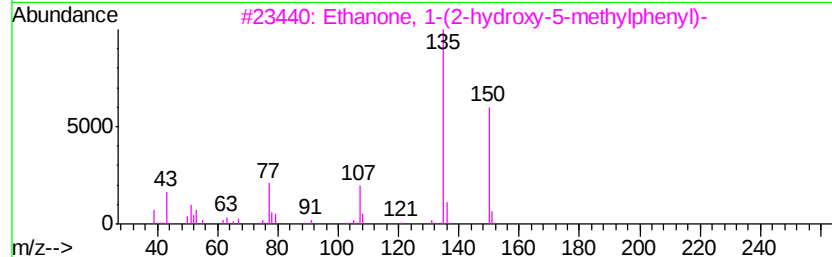
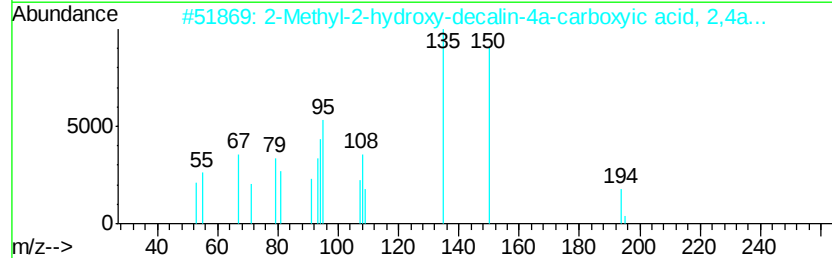
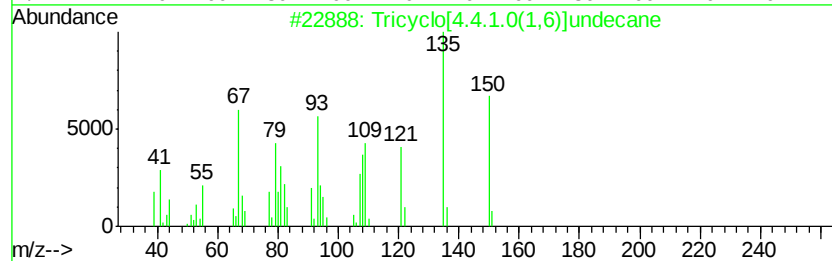
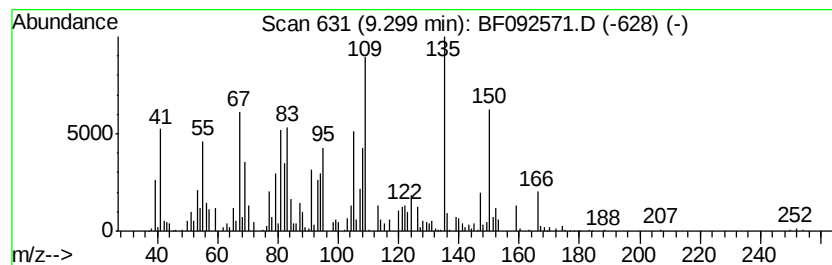
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown9.30 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.30	22.07 ng	2415750	Acenaphthene-d10	10.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[4.4.1.0(1,6)]undecane	150	C11H18	006571-73-9	45
2		2-Methyl-2-hydroxy-decalin-4a-ca...	194	C12H18O2	1000146-22-6	43
3		Ethanone, 1-(2-hydroxy-5-methylp...	150	C9H10O2	001450-72-2	38
4		Ethanone, 1-(6,6-dimethylbicyclo...	150	C10H14O	024555-40-6	35
5		2,5-Dimethyl-3-isopropylpyrazine	150	C9H14N2	013610-20-3	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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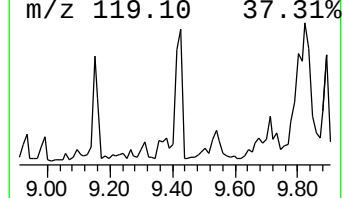
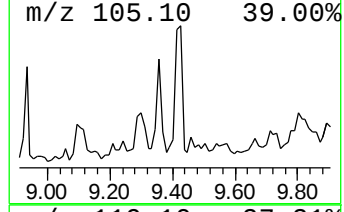
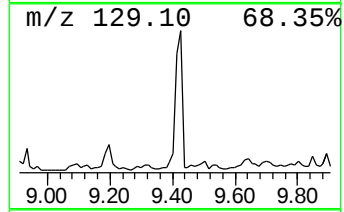
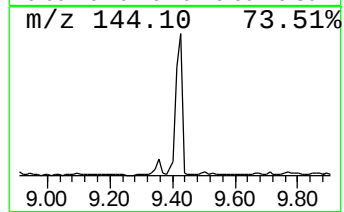
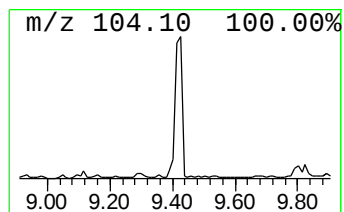
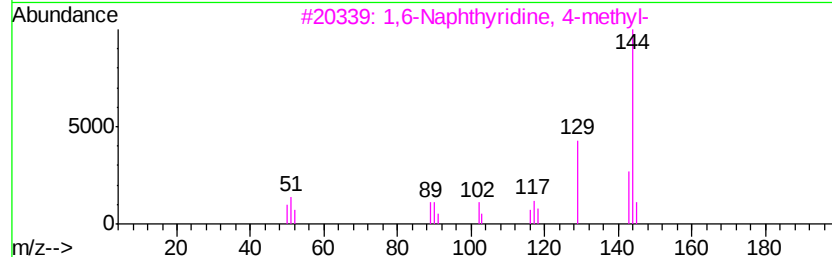
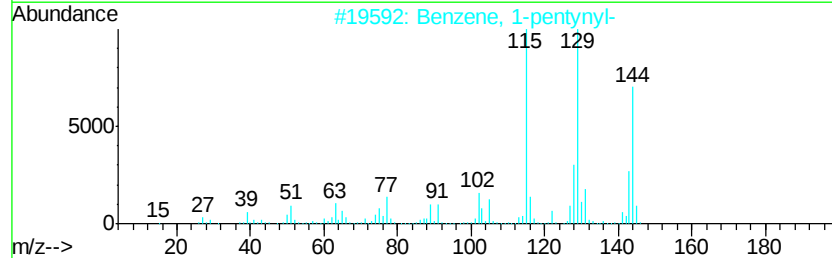
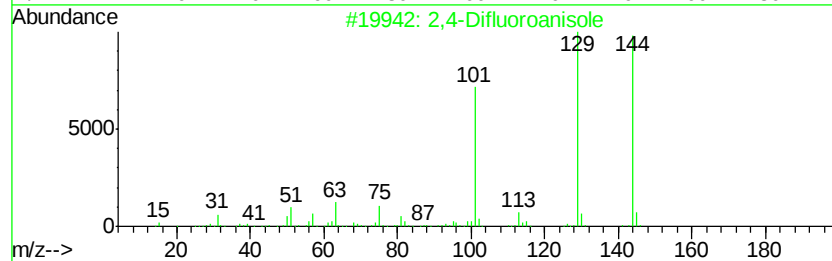
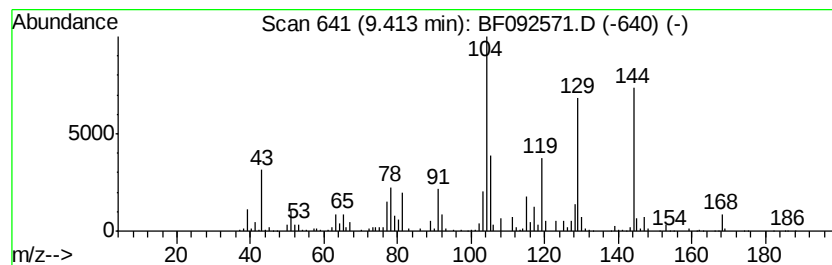
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown9.41 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.41	22.89 ng	2506400	Acenaphthene-d10	10.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Difluoroanisole	144	C7H6F2O	000452-10-8	38
2		Benzene, 1-pentynyl-	144	C11H12	004250-81-1	35
3		1,6-Naphthvridine, 4-methyl-	144	C9H8N2	007675-30-1	35
4		1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	27
5		2-Azetidinone, 4-phenyl-	147	C9H9NO	005661-55-2	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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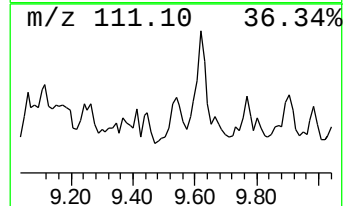
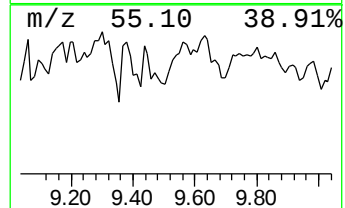
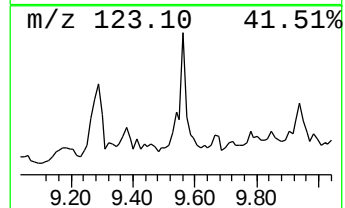
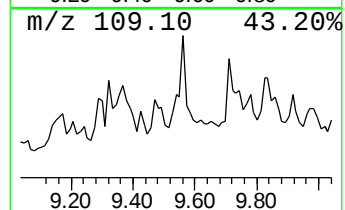
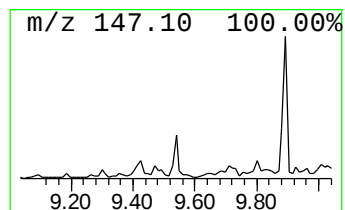
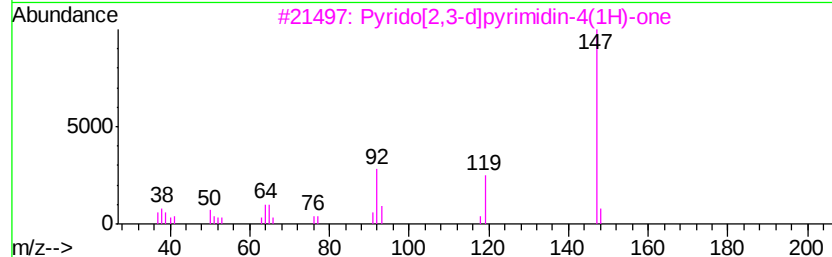
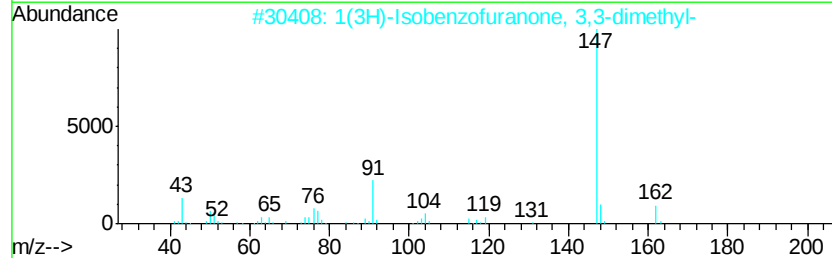
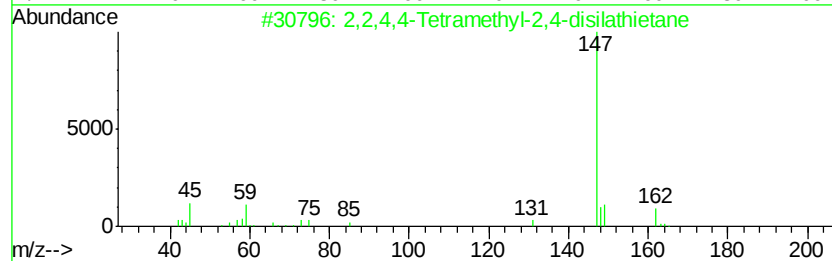
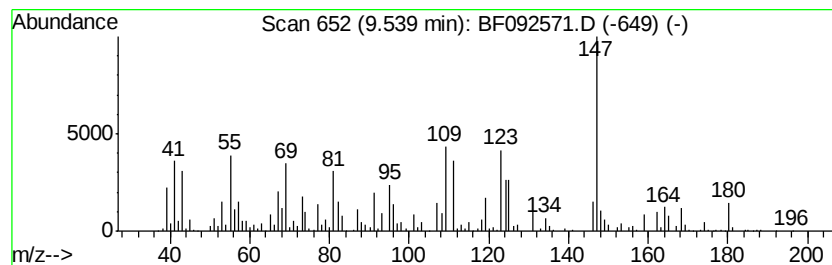
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown9.54 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.54	14.48 ng	1584830	Acenaphthene-d10	10.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,4,4-Tetramethyl-2,4-disilath...	162	C5H14SSi2	091521-58-3	18		
2	1(3H)-Isobenzofuranone, 3,3-dime...	162	C10H10O2	001689-09-4	18		
3	Pyrrolo[2,3-d]pyrimidin-4(1H)-one	147	C7H5N3O	024410-19-3	18		
4	Ethanone, 1,1'-(1,3-phenylene)bis-	162	C10H10O2	006781-42-6	18		
5	o-Diacetylbenzene	162	C10H10O2	000704-00-7	18		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

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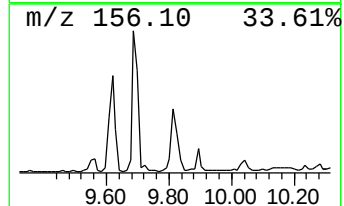
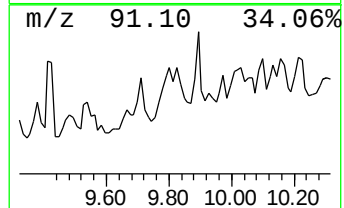
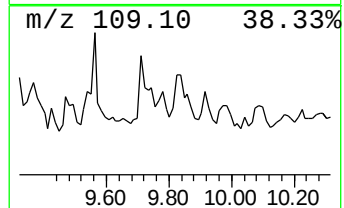
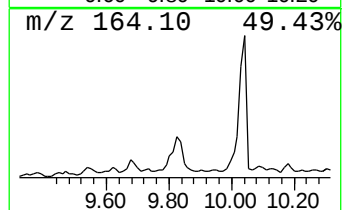
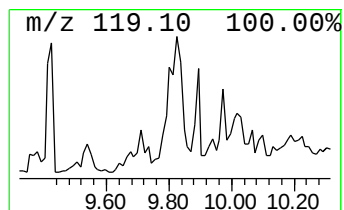
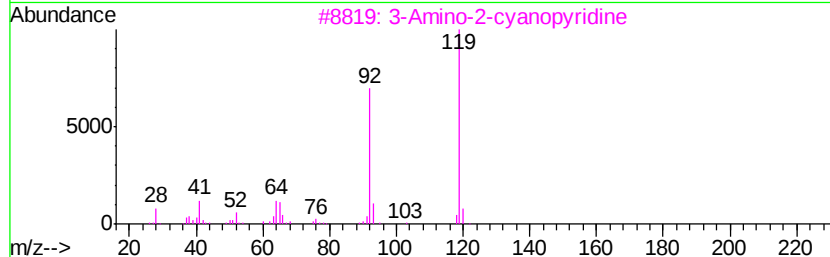
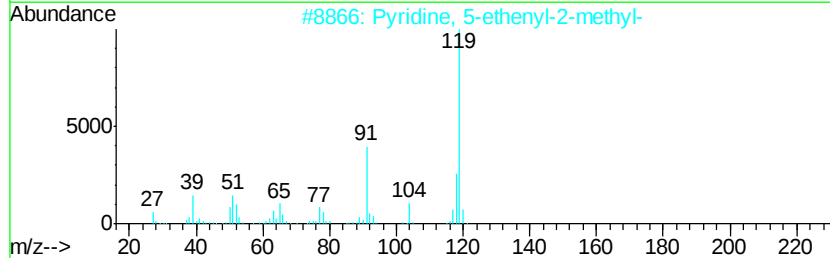
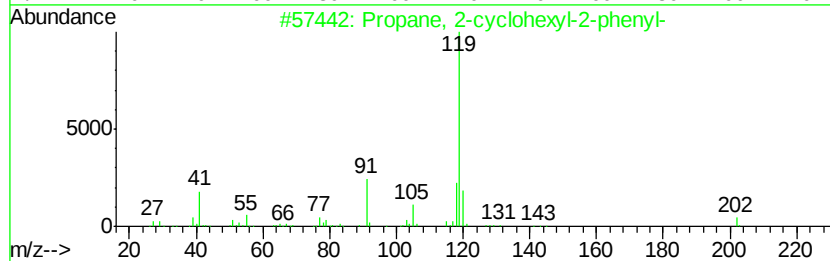
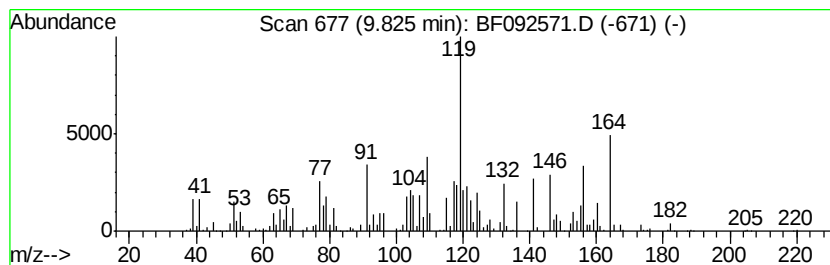
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown9.82 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	34.38 ng	3764460	Acenaphthene-d10	10.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2-cyclohexyl-2-phenyl-	202	C15H22	025683-97-0	27
2		Pyridine, 5-ethenyl-2-methyl-	119	C8H9N	000140-76-1	25
3		3-Amino-2-cyanopyridine	119	C6H5N3	042242-11-5	14
4		Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	14
5		.alpha.-Bromomesitylene	198	C9H11Br	027129-86-8	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
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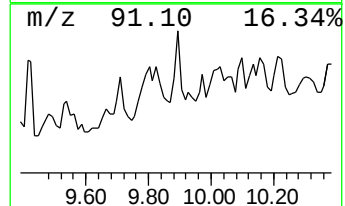
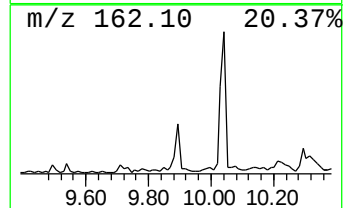
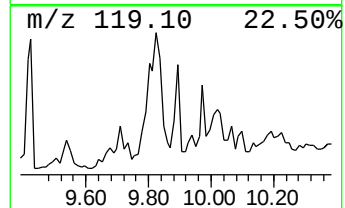
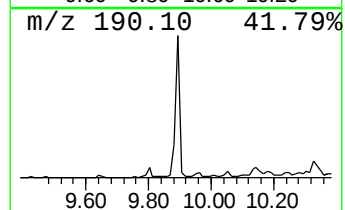
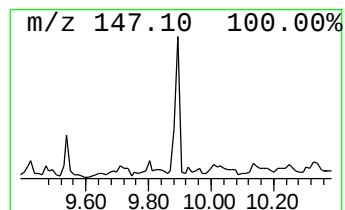
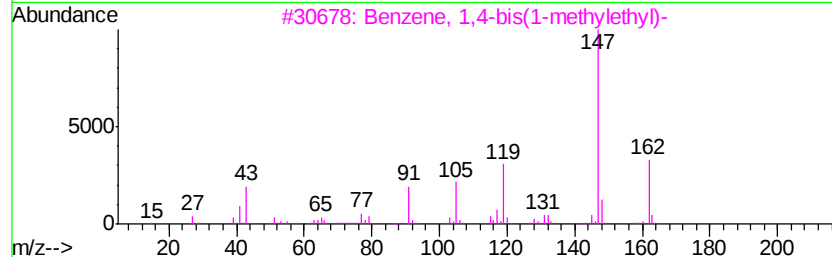
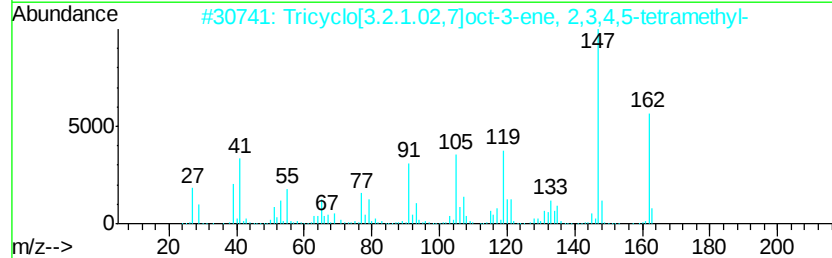
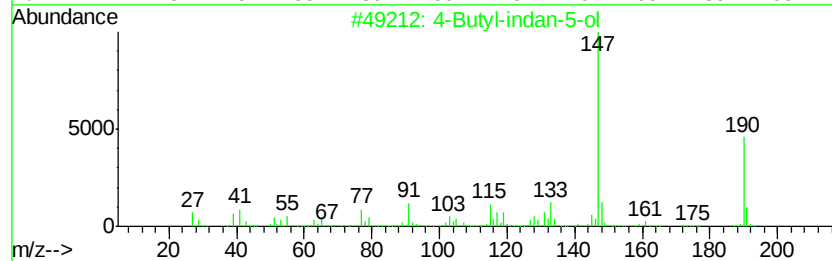
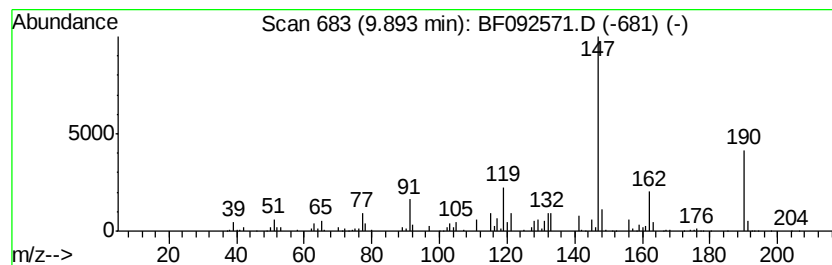
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 4-Butyl-indan-5-ol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	12.46 ng	1364570	Acenaphthene-d10	10.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4-Butyl-indan-5-ol	190 C13H18O	1000194-67-3 70
2		Tricyclo[3.2.1.0 ^{2,7}]oct-3-ene, 2...	162 C12H18	062338-44-7 68
3		Benzene, 1,4-bis(1-methylethyl)-	162 C12H18	000100-18-5 64
4		Benzene, 1-(1,1-dimethylethyl)-3...	162 C12H18	000098-19-1 60
5		Benzene, 1-(1,1-dimethylethyl)-3...	162 C12H18	014411-56-4 60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
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Misc :
ALS Vial : 56 Sample Multiplier: 1

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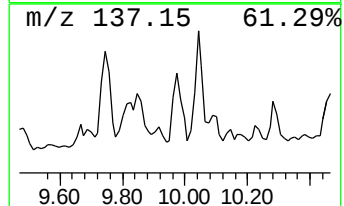
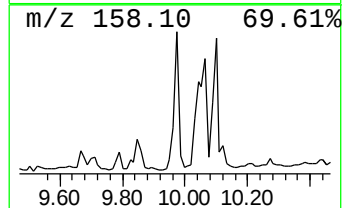
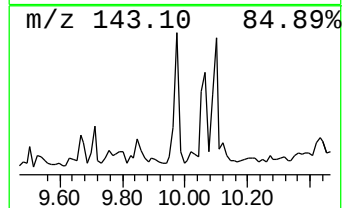
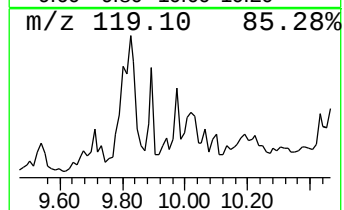
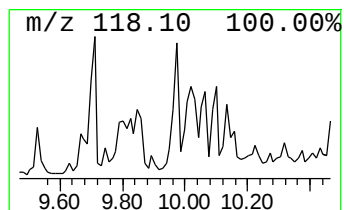
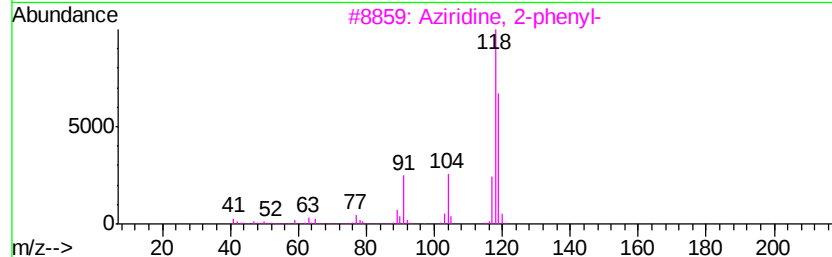
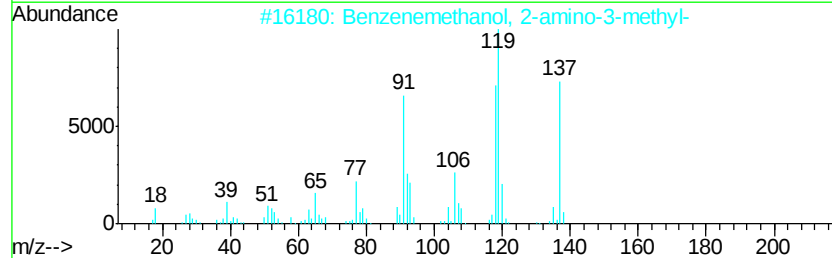
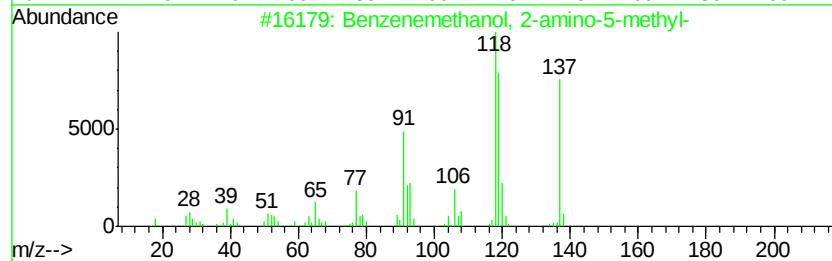
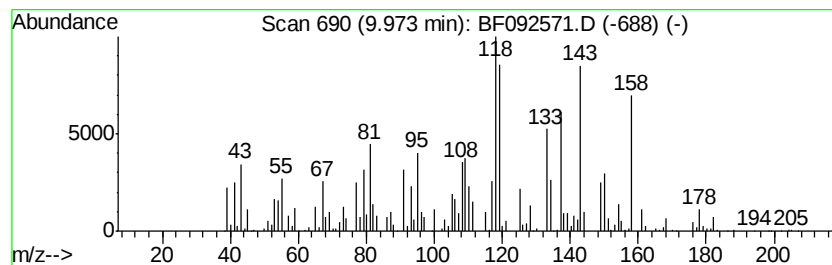
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown9.97 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.97	14.19 ng	1553410	Acenaphthene-d10	10.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanol, 2-amino-5-methyl-	137	C8H11NO	034897-84-2	42
2		Benzenemethanol, 2-amino-3-methyl-	137	C8H11NO	057772-50-6	25
3		Aziridine, 2-phenyl-	119	C8H9N	001499-00-9	22
4		1H-Indole, 2,3-dihydro-	119	C8H9N	000496-15-1	22
5		3-Cyclohexene-1-ethanol, .alpha....	220	C15H24O	055780-93-3	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-02

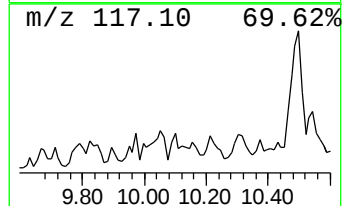
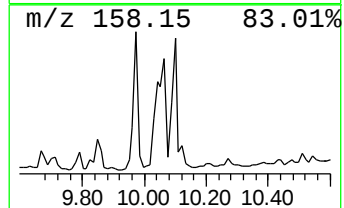
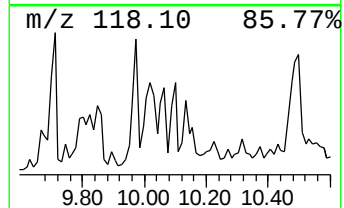
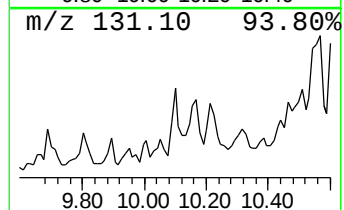
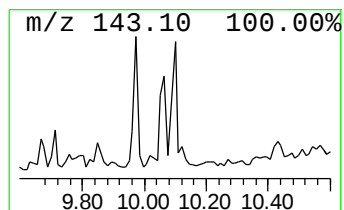
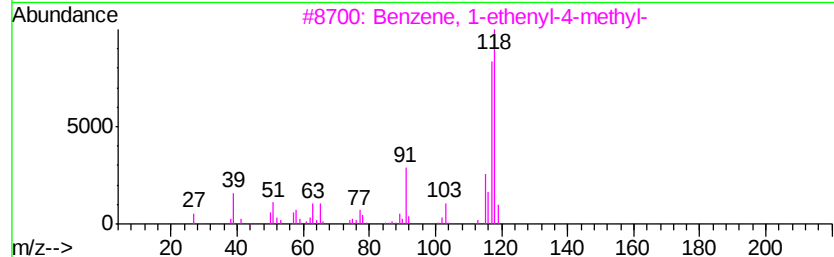
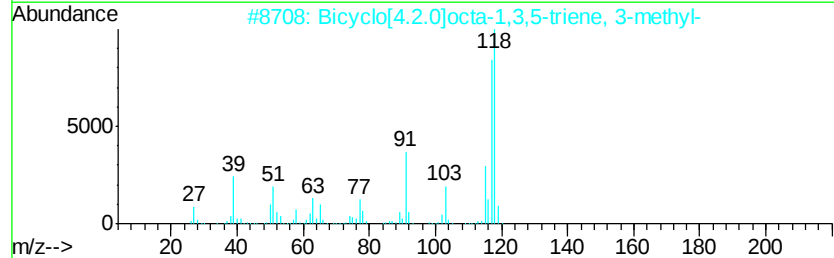
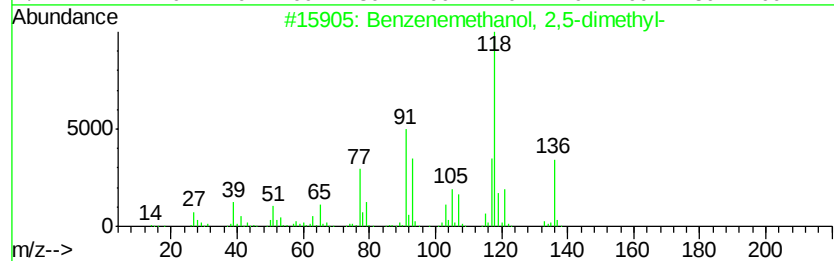
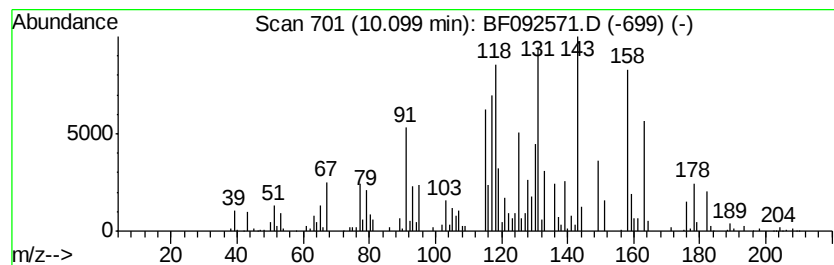
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown10.10 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.10	8.70 ng	952711	Acenaphthene-d10	10.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanol, 2,5-dimethyl-	136	C9H12O	053957-33-8	25
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	15
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	15
4		Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	15
5		Cyclohexene, 1-phenyl-	158	C12H14	000771-98-2	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
Data File : BF092571.D
Acq On : 27 Jan 2017 12:27
Operator : SJ/MA
Sample : I1353-07
Misc :
ALS Vial : 56 Sample Multiplier: 1

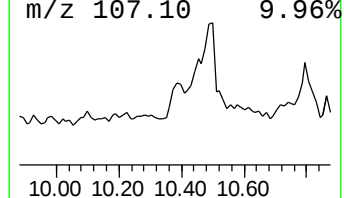
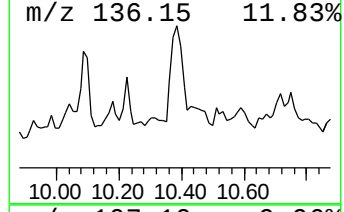
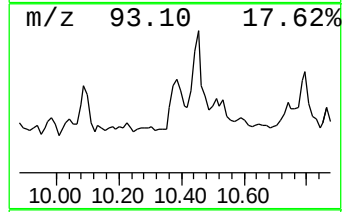
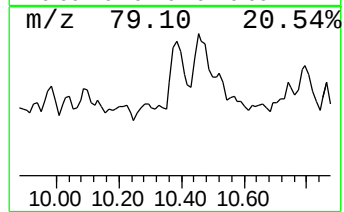
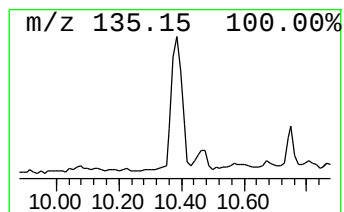
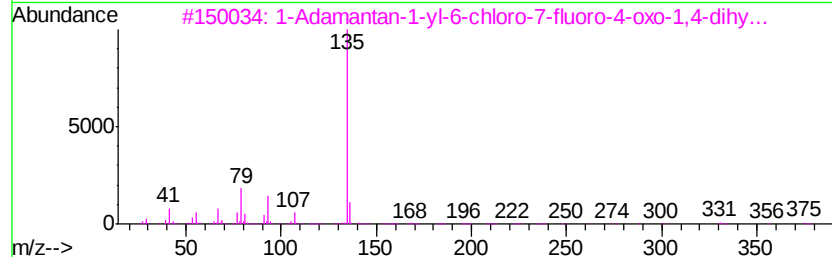
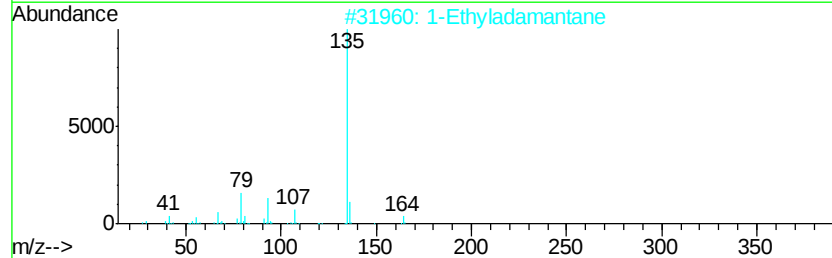
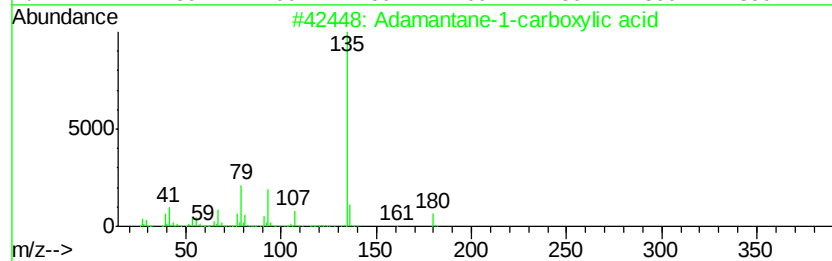
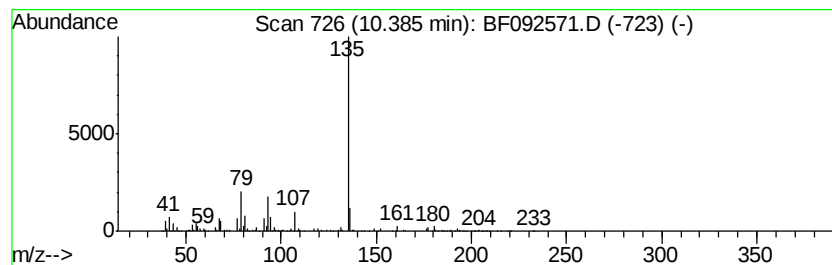
Instrument :
BNA_F
ClientSampleId :
W-02

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Adamantane-1-carboxylic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	17.82 ng	1951400	Acenaphthene-d10	10.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Adamantane-1-carboxylic acid	180 C11H16O2	000828-51-3 90
2		1-Ethyladamantane	164 C12H20	000770-69-4 87
3		1-Adamantan-1-yl-6-chloro-7-fluo...	375 C20H19ClFN03	1000210-85-1 86
4		1-Adamantanecarboxylic acid chloride	198 C11H15ClO	002094-72-6 86
5		1-Adamantyl bromomethyl ketone	256 C12H17BrO	005122-82-7 86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012617\
 Data File : BF092571.D
 Acq On : 27 Jan 2017 12:27
 Operator : SJ/MA
 Sample : I1353-07
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-02

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012417.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-Pentanone, 4-hy...	5.16	12.9	ng	708478	1	6.97	1094020	20.0
unknown6.75	6.75	76.6	ng	4190440	1	6.97	1094020	20.0
Benzene, 1,3-diet...	7.22	14.2	ng	778423	1	6.97	1094020	20.0
Benzene, 1,2-diet...	7.31	26.6	ng	1456920	1	6.97	1094020	20.0
unknown7.41	7.41	14.7	ng	804685	1	6.97	1094020	20.0
Benzene, 1-ethyl-...	7.62	14.6	ng	798975	1	6.97	1094020	20.0
1H-Indene, 2,3-di...	7.94	7.3	ng	2096640	2	8.27	5709650	20.0
Naphthalene, 1,2,...	8.46	9.8	ng	2808740	2	8.27	5709650	20.0
unknown8.73	8.73	9.8	ng	2788250	2	8.27	5709650	20.0
Naphthalene, 1,2,...	8.93	11.4	ng	3263560	2	8.27	5709650	20.0
unknown9.16	9.16	17.5	ng	1920500	3	10.04	2189620	20.0
unknown9.30	9.30	22.1	ng	2415750	3	10.04	2189620	20.0
unknown9.41	9.41	22.9	ng	2506400	3	10.04	2189620	20.0
unknown9.54	9.54	14.5	ng	1584830	3	10.04	2189620	20.0
unknown9.82	9.82	34.4	ng	3764460	3	10.04	2189620	20.0
4-Butyl-indan-5-ol	9.89	12.5	ng	1364570	3	10.04	2189620	20.0
unknown9.97	9.97	14.2	ng	1553410	3	10.04	2189620	20.0
unknown10.10	10.10	8.7	ng	952711	3	10.04	2189620	20.0
Adamantane-1-carb...	10.38	17.8	ng	1951400	3	10.04	2189620	20.0