

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
 Data File : BF084557.D
 Acq On : 19 Jan 2016 17:34
 Operator : SJ/UM
 Sample : H1169-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-004(0-2)

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-BF123115.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	3.699	131	141	144	rBV2	73293	184909	3.13%	0.190%
2	4.887	240	245	249	rBV2	119783	244521	4.14%	0.252%
3	4.967	249	252	255	rBV	1881585	3369003	57.00%	3.471%
4	5.173	267	270	271	rBV	180416	223103	3.77%	0.230%
5	5.264	275	278	283	rVB3	95513	190225	3.22%	0.196%
6	5.367	283	287	288	rBV2	257142	469143	7.94%	0.483%
7	5.402	288	290	292	rVV	5478778	5378269	90.99%	5.541%
8	5.630	308	310	313	rVB3	106561	190389	3.22%	0.196%
9	5.687	313	315	318	rVB	223051	214770	3.63%	0.221%
10	5.802	322	325	327	rBV2	280124	427947	7.24%	0.441%
11	5.859	327	330	334	rBV3	106432	260028	4.40%	0.268%
12	5.939	334	337	339	rVB2	260800	336010	5.68%	0.346%
13	5.984	339	341	344	rBV2	108570	212265	3.59%	0.219%
14	6.030	344	345	348	rVB	443292	554088	9.37%	0.571%
15	6.087	348	350	352	rBV	203257	196512	3.32%	0.202%
16	6.224	359	362	364	rBV	247295	348853	5.90%	0.359%
17	6.293	364	368	369	rVV2	190556	255414	4.32%	0.263%
18	6.316	369	370	374	rVB3	229555	359050	6.07%	0.370%
19	6.385	374	376	378	rBV	223224	228901	3.87%	0.236%
20	6.442	378	381	383	rBV	5125782	5910519	100.00%	6.089%
21	6.567	389	392	394	rVB	4693608	5726431	96.89%	5.899%
22	6.636	394	398	400	rVB3	348806	623796	10.55%	0.643%
23	6.796	410	412	415	rVB2	1208533	1754494	29.68%	1.807%
24	6.853	415	417	419	rBV	311328	345425	5.84%	0.356%
25	6.945	423	425	427	rVV	3593563	4069828	68.86%	4.193%
26	7.070	429	436	438	rBV3	233376	633320	10.72%	0.652%
27	7.116	438	440	441	rBV2	434508	464928	7.87%	0.479%
28	7.139	441	442	444	rVB	200923	181251	3.07%	0.187%
29	7.185	444	446	448	rBV	381703	401732	6.80%	0.414%
30	7.287	451	455	456	rBV	210289	300029	5.08%	0.309%
31	7.356	456	461	463	rBV2	2854142	4432973	75.00%	4.567%
32	7.402	463	465	466	rVB	485055	432555	7.32%	0.446%
33	7.447	466	469	470	rBV3	254583	361702	6.12%	0.373%
34	7.482	470	472	473	rBV	222672	243144	4.11%	0.250%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	7.516	473	475	477 rVB	181949	231894	3.92%	0.239%
36	7.573	477	480	481 rBV	548385	890380	15.06%	0.917%
37	7.596	481	482	484 rVB	837416	694721	11.75%	0.716%
38	7.733	487	494	495 rBV6	386194	1258858	21.30%	1.297%
39	7.768	495	497	499 rVB2	264494	351411	5.95%	0.362%
40	7.813	499	501	506 rBV4	761886	1491494	25.23%	1.537%
41	7.905	506	509	511 rVB4	290936	570021	9.64%	0.587%
42	7.950	511	513	515 rBV	261324	268542	4.54%	0.277%
43	8.042	518	521	522 rBV3	347823	672348	11.38%	0.693%
44	8.076	522	524	525 rVV2	2737399	3206810	54.26%	3.304%
45	8.145	528	530	537 rVB3	521735	1436261	24.30%	1.480%
46	8.248	537	539	541 rBV3	147243	292406	4.95%	0.301%
47	8.373	546	550	554 rVV6	459759	1409671	23.85%	1.452%
48	8.465	556	558	560 rVV	835654	763135	12.91%	0.786%
49	8.510	560	562	564 rVV	1037799	1201315	20.33%	1.238%
50	8.545	564	565	567 rVV	516718	494287	8.36%	0.509%
51	8.590	567	569	570 rVV	213503	321353	5.44%	0.331%
52	8.625	570	572	575 rVV3	149552	362566	6.13%	0.374%
53	8.682	575	577	578 rVV2	986291	1133621	19.18%	1.168%
54	8.716	578	580	583 rVV2	326341	707089	11.96%	0.728%
55	8.796	583	587	588 rVB2	366410	799722	13.53%	0.824%
56	8.830	588	590	591 rBV	123323	178936	3.03%	0.184%
57	8.888	593	595	597 rVV	395210	487988	8.26%	0.503%
58	8.990	600	604	606 rVV3	218732	501547	8.49%	0.517%
59	9.036	606	608	610 rVV3	122412	246940	4.18%	0.254%
60	9.070	610	611	613 rVV2	177017	235448	3.98%	0.243%
61	9.105	613	614	616 rVV	444014	496564	8.40%	0.512%
62	9.162	616	619	621 rVB	5998859	5874987	99.40%	6.052%
63	9.242	623	626	628 rBV2	671184	910245	15.40%	0.938%
64	9.299	628	631	634 rVV5	152489	476884	8.07%	0.491%
65	9.413	638	641	644 rVV3	247384	511311	8.65%	0.527%
66	9.493	646	648	651 rVV2	498517	750729	12.70%	0.773%
67	9.562	651	654	656 rVV3	1125705	2386674	40.38%	2.459%
68	9.596	656	657	661 rVB2	235669	353184	5.98%	0.364%
69	9.699	663	666	668 rBV3	388368	525612	8.89%	0.541%
70	9.745	668	670	671 rVB	815202	624879	10.57%	0.644%
71	9.768	671	672	674 rBV	295295	423763	7.17%	0.437%

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72	9.836	674	678	680	rVV	1961710	2219711	37.56%	2.287%
73	9.939	686	687	689	rVB2	180783	220115	3.72%	0.227%
74	9.996	689	692	694	rBV3	105154	246262	4.17%	0.254%
75	10.031	694	695	697	rVB	273289	265565	4.49%	0.274%
76	10.156	703	706	709	rBV4	199818	443484	7.50%	0.457%
77	10.236	712	713	715	rBV	273512	319050	5.40%	0.329%
78	10.271	715	716	717	rVB	476014	327905	5.55%	0.338%
79	10.488	733	735	738	rVB	483519	614733	10.40%	0.633%
80	10.625	744	747	749	rBV	3274944	3701581	62.63%	3.813%
81	10.762	756	759	761	rBV	981982	1695841	28.69%	1.747%
82	11.196	795	797	798	rBV	275530	353874	5.99%	0.365%
83	11.219	798	799	802	rVB	434946	449035	7.60%	0.463%
84	11.311	805	807	811	rBV	1560857	1907177	32.27%	1.965%
85	11.619	832	834	835	rVB	283647	219466	3.71%	0.226%
86	11.928	859	861	866	rVB4	157320	336230	5.69%	0.346%
87	12.019	868	869	874	rVB	149794	192539	3.26%	0.198%
88	12.408	901	903	905	rVB2	151776	207972	3.52%	0.214%
89	12.522	911	913	916	rBV	751659	654399	11.07%	0.674%
90	12.751	930	933	935	rBV	678188	668530	11.31%	0.689%
91	12.899	944	946	949	rVB	3883845	3939283	66.65%	4.058%
92	13.814	1023	1026	1035	rBV3	183171	416341	7.04%	0.429%
93	13.951	1035	1038	1039	rBV2	1275185	1575535	26.66%	1.623%
94	14.957	1123	1126	1131	rBV	340865	624838	10.57%	0.644%
95	15.243	1148	1151	1154	rVV	171904	219284	3.71%	0.226%
96	15.300	1154	1156	1158	rVV	197939	259000	4.38%	0.267%
97	15.357	1158	1161	1167	rVB	945584	1251626	21.18%	1.289%
98	16.706	1277	1279	1283	rBV2	112125	201008	3.40%	0.207%
99	17.117	1311	1315	1322	rVB	99419	240435	4.07%	0.248%
100	19.151	1491	1493	1502	rVB6	60546	230914	3.91%	0.238%

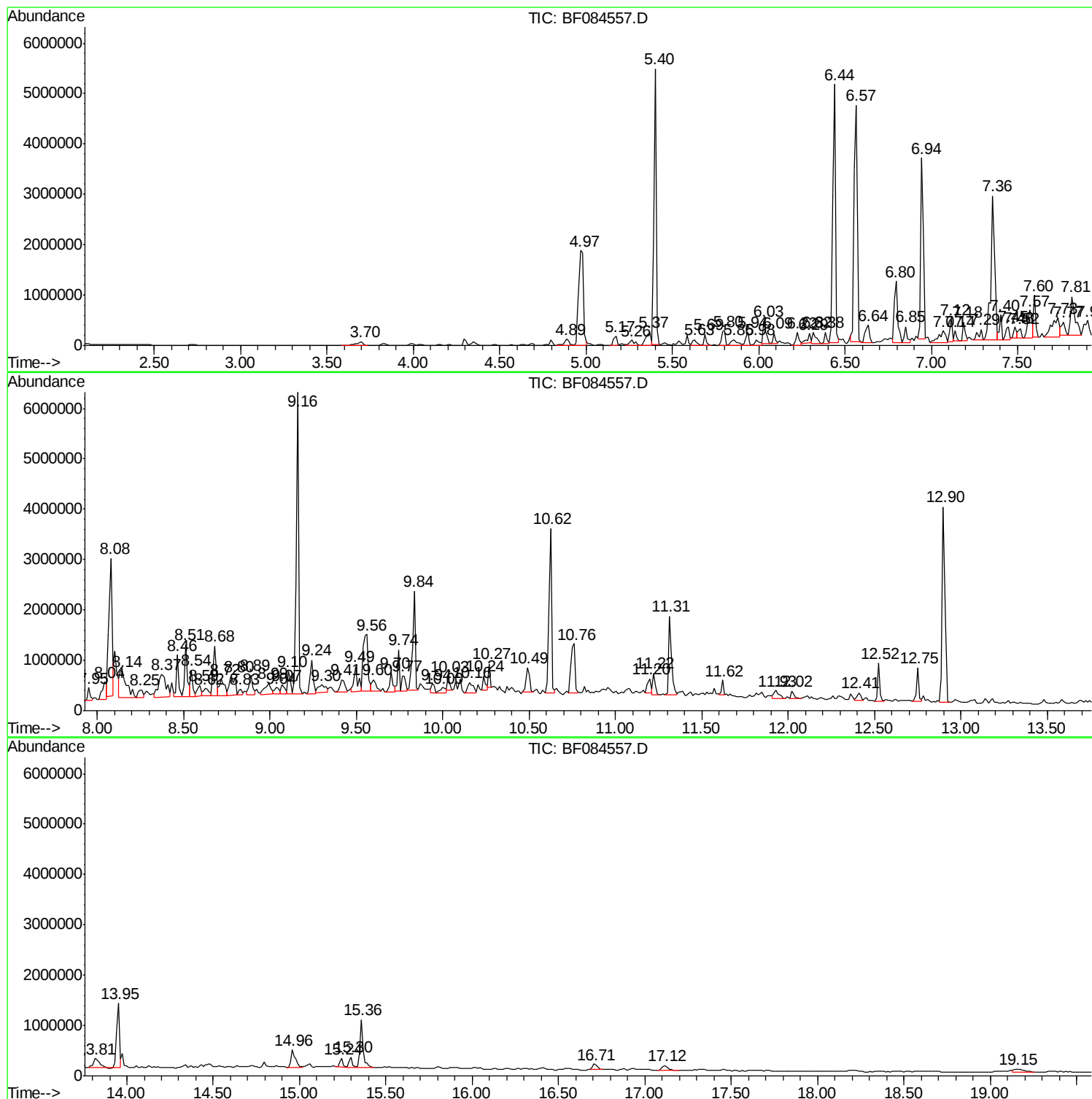
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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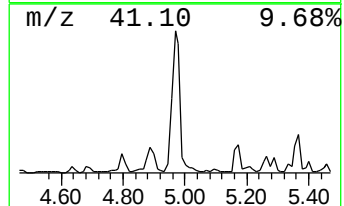
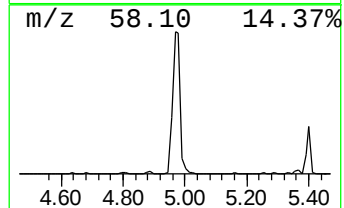
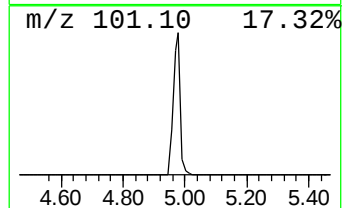
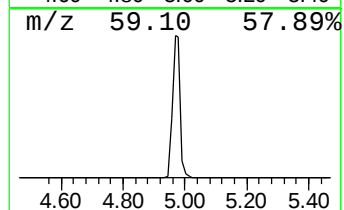
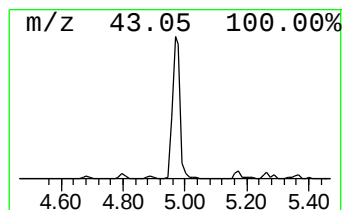
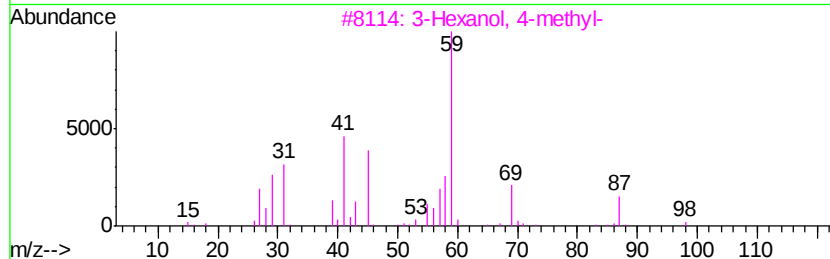
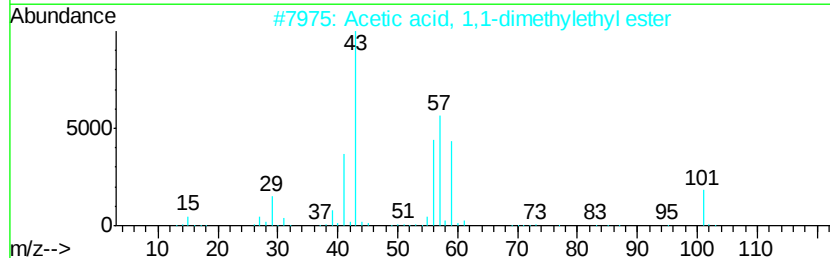
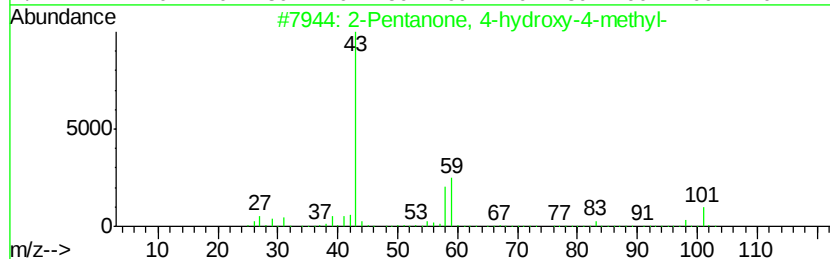
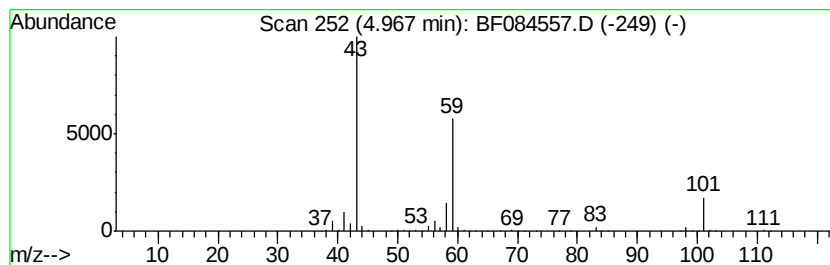
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	38.40 ng	3369000	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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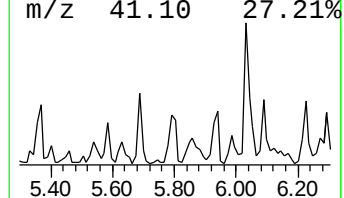
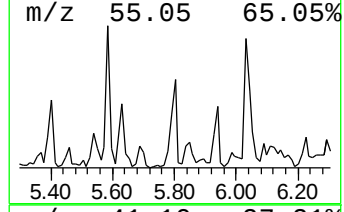
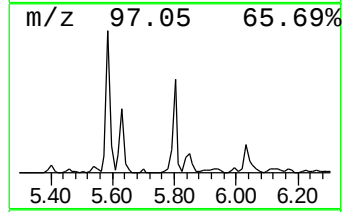
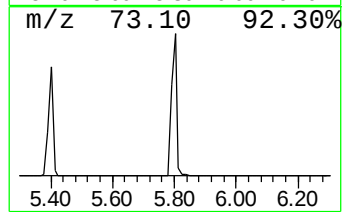
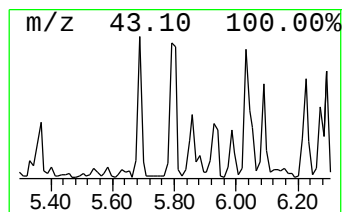
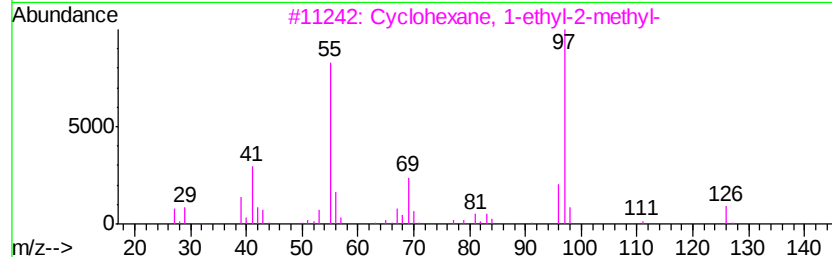
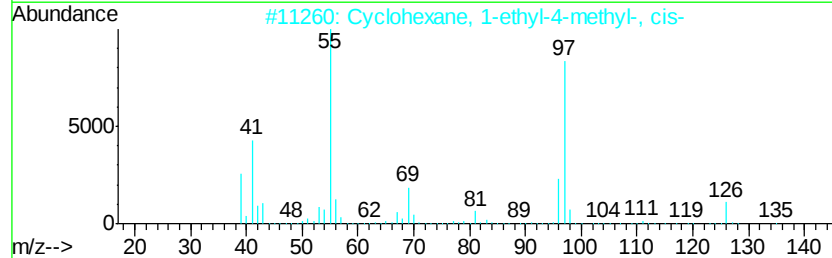
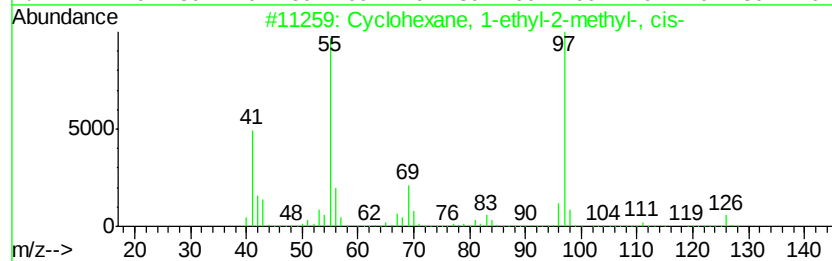
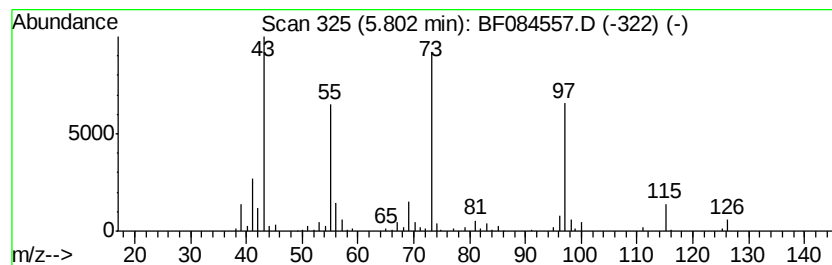
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.80	4.88 ng	427947	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	70
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	58
3		Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	58
4		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	52
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	52



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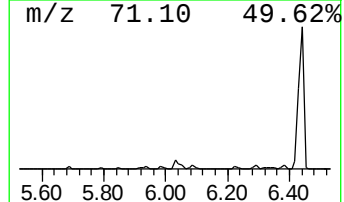
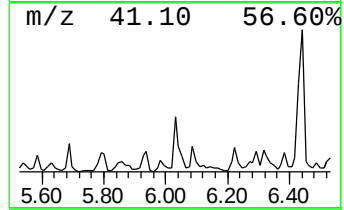
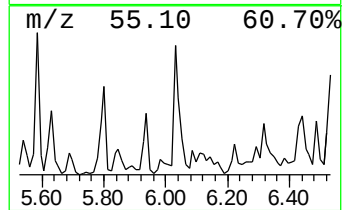
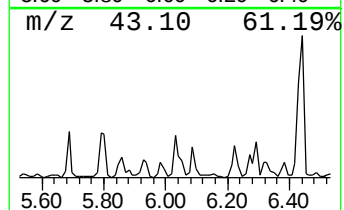
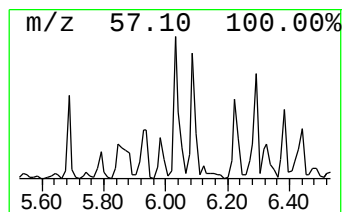
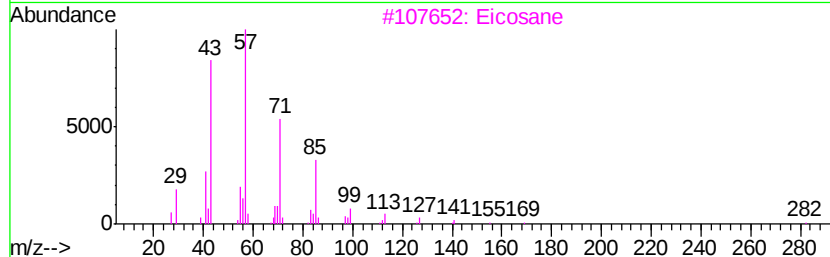
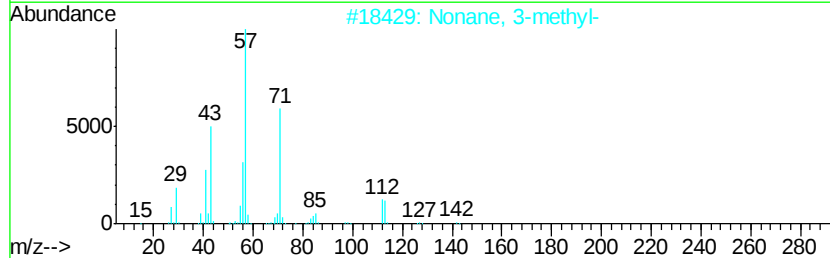
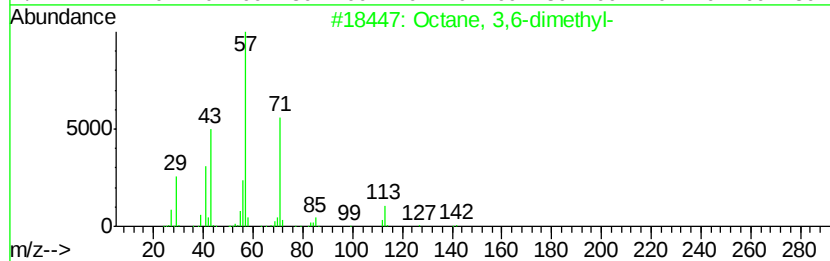
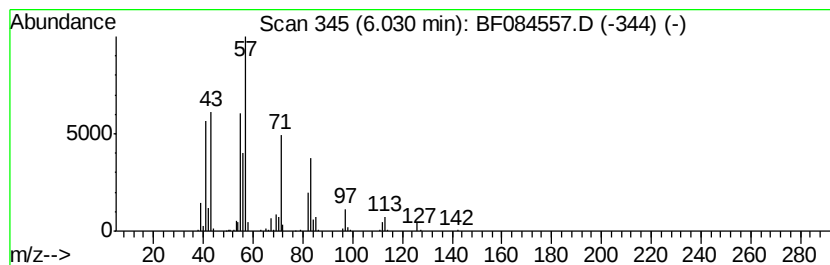
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TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.03 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.03	6.32 ng	554088	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	46
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	43
3		Eicosane	282	C20H42	000112-95-8	38
4		Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	35
5		Tetracosane	338	C24H50	000646-31-1	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
Sample : H1169-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-004(0-2)

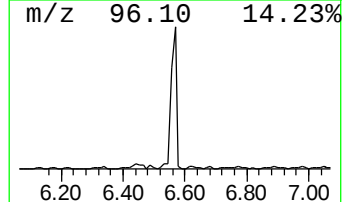
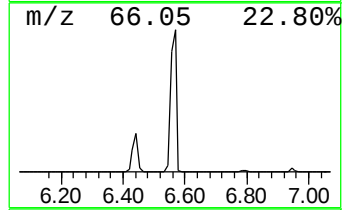
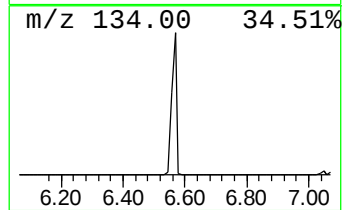
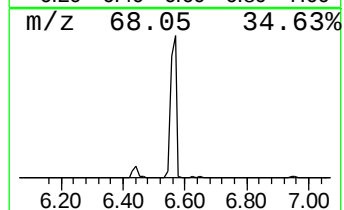
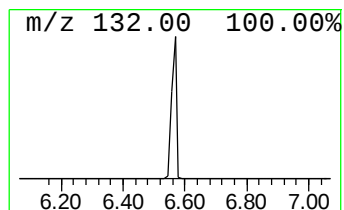
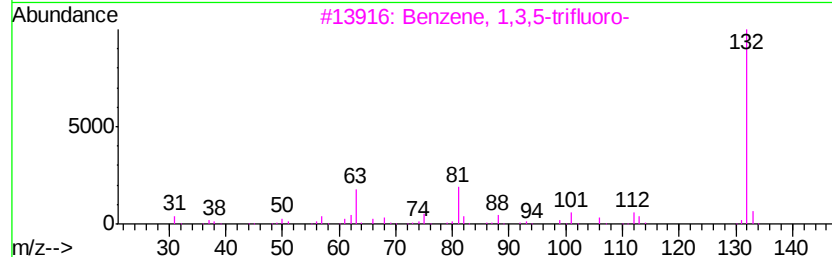
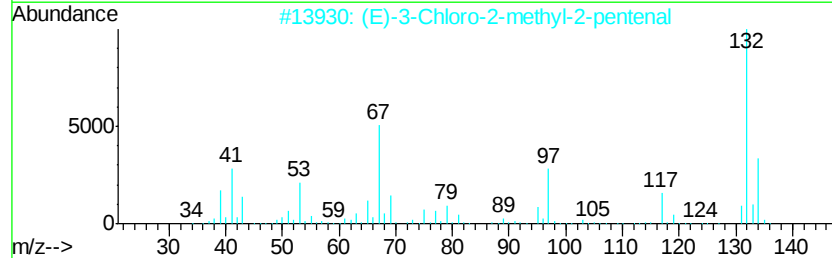
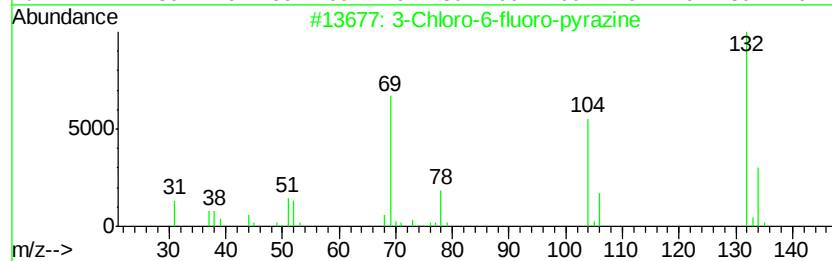
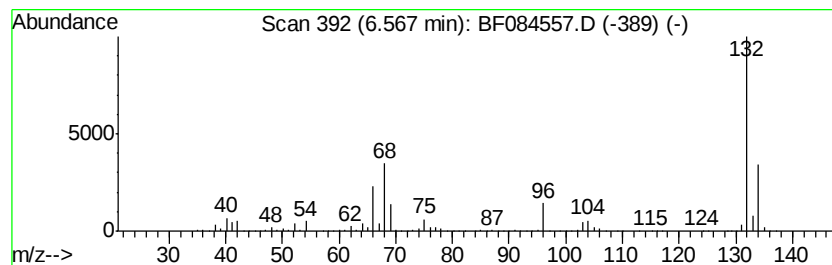
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	65.28 ng	5726430	1,4-Dichlorobenzene-d4	6.80

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	35
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
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Misc :
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Instrument :
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ClientSampleId :
SP-004(0-2)

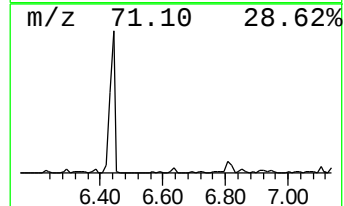
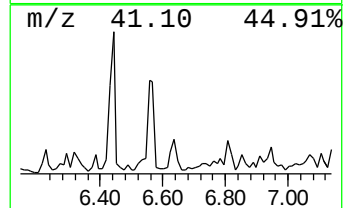
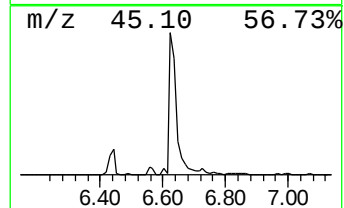
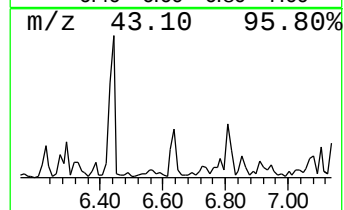
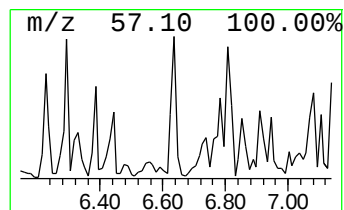
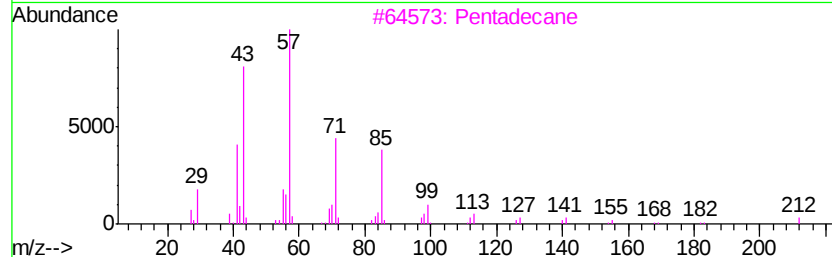
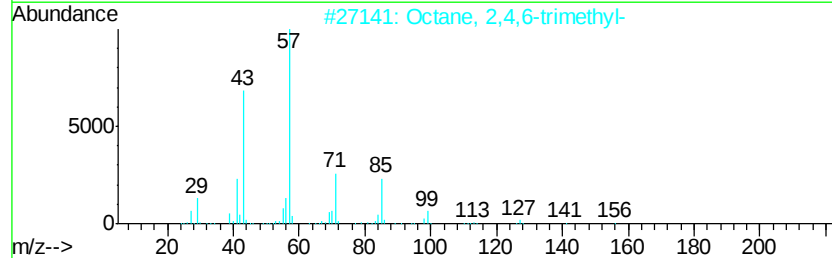
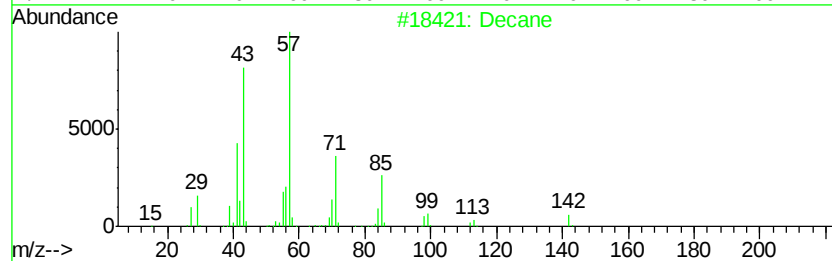
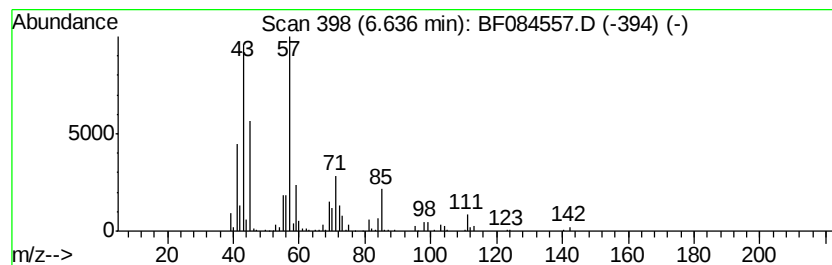
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Decane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	7.11 ng	623796	1,4-Dichlorobenzene-d4	6.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	50
2	Octane, 2,4,6-trimethyl-		156	C11H24	062016-37-9	38
3	Pentadecane		212	C15H32	000629-62-9	35
4	Ethanol, 2-(2-ethoxyethoxy)-		134	C6H14O3	000111-90-0	35
5	Dotriacontane		451	C32H66	000544-85-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
Sample : H1169-07
Misc :
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ClientSampleId :
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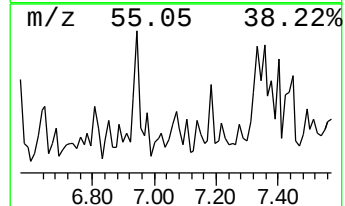
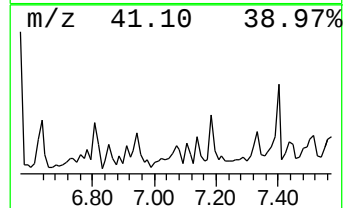
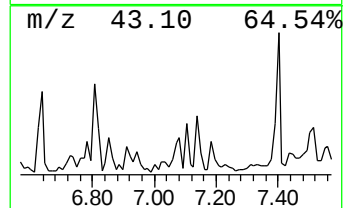
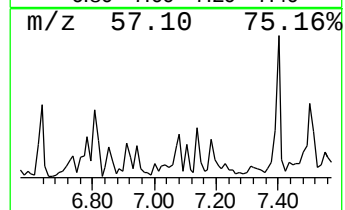
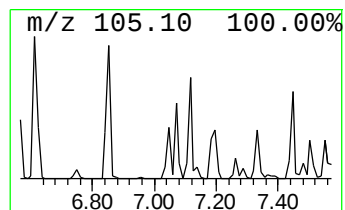
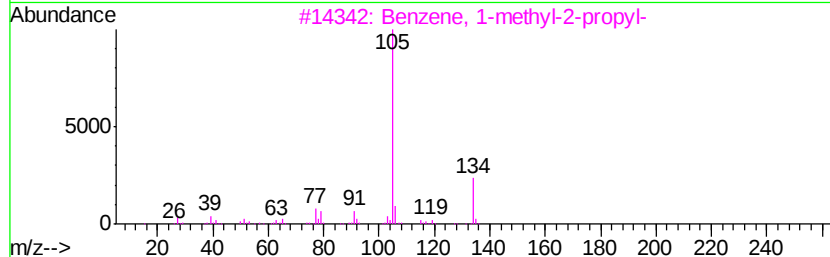
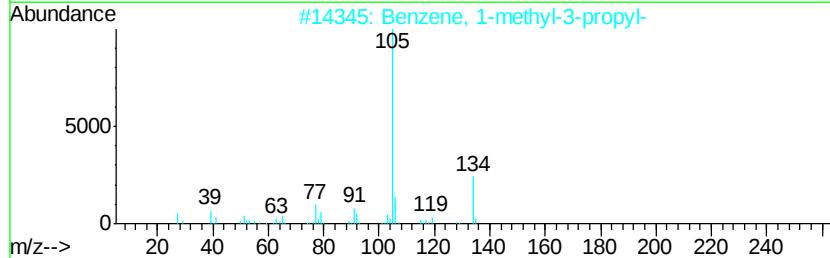
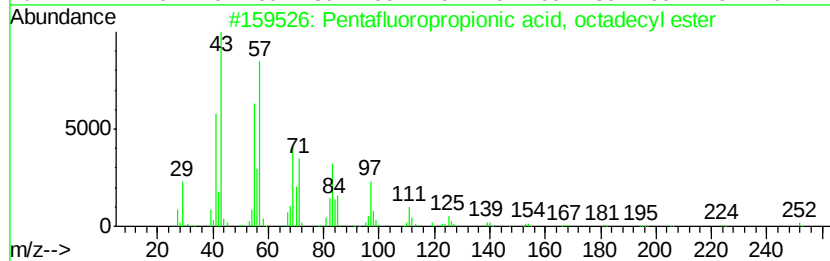
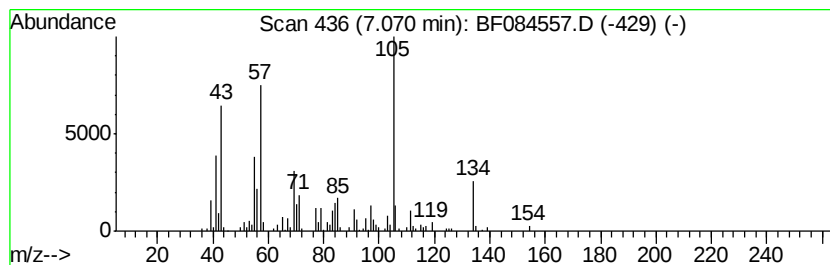
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown7.07 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	7.22 ng	633320	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	1000280-07-7	43
2		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	42
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	42
4		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	42
5		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
Sample : H1169-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

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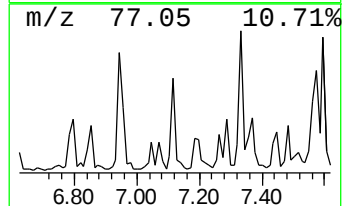
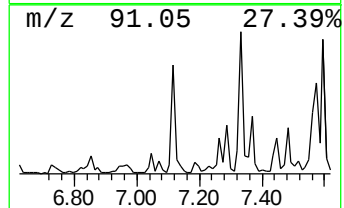
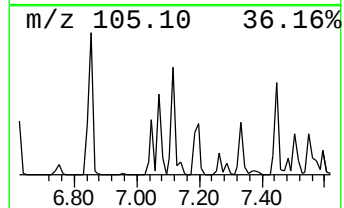
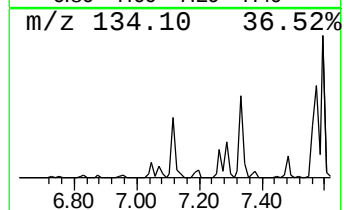
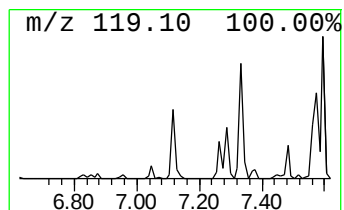
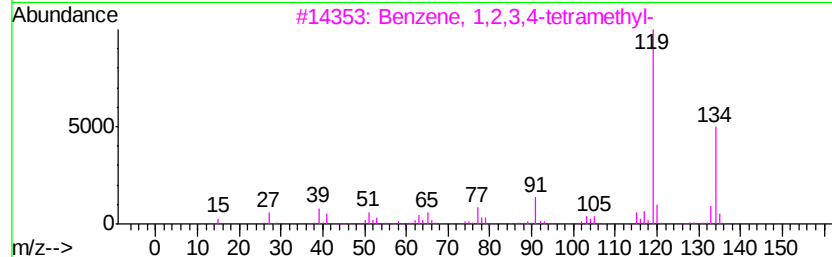
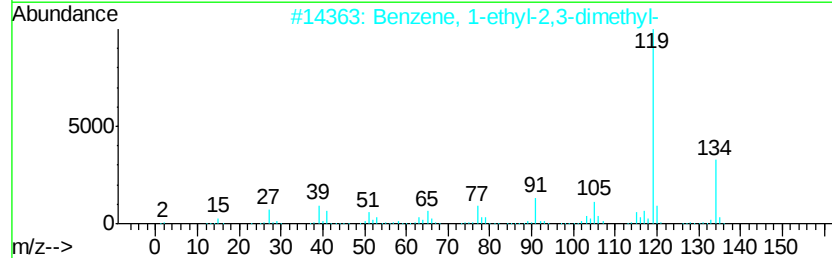
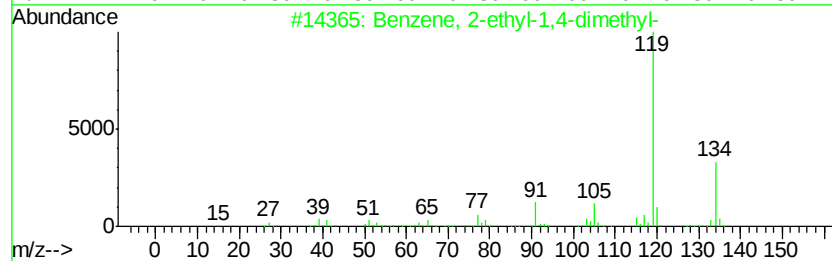
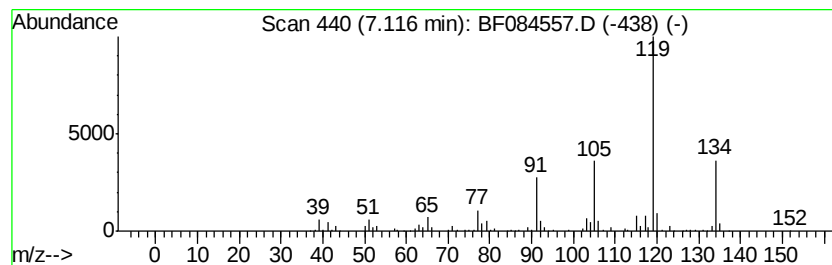
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	5.30 ng	464928	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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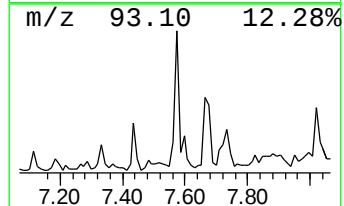
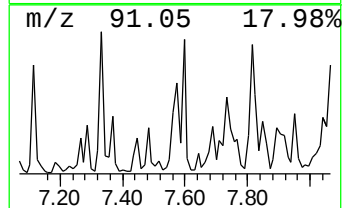
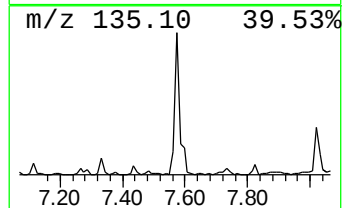
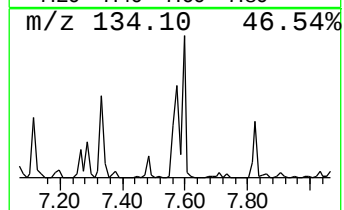
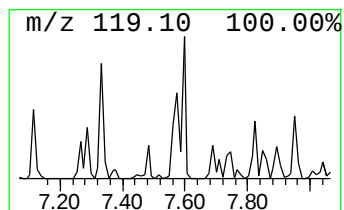
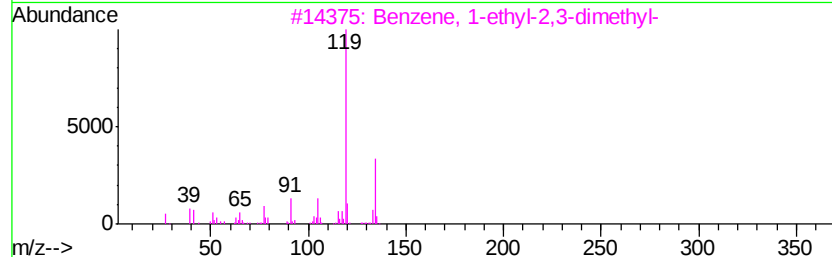
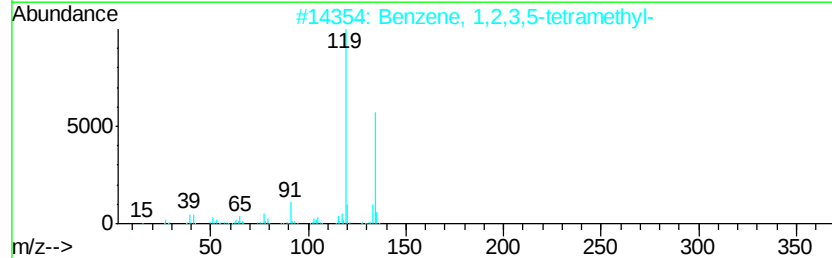
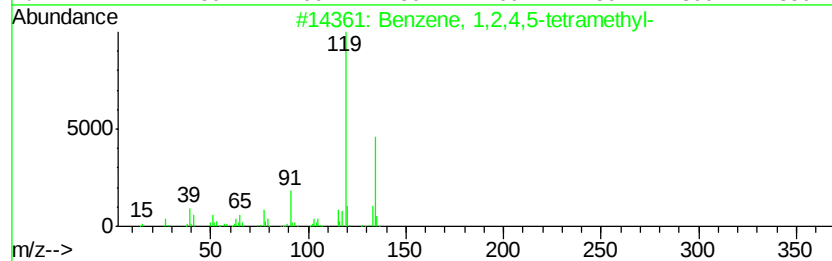
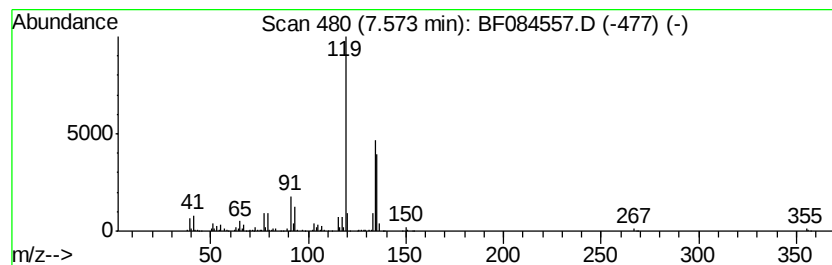
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	5.55 ng	890380	Naphthalene-d8	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	93
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	76
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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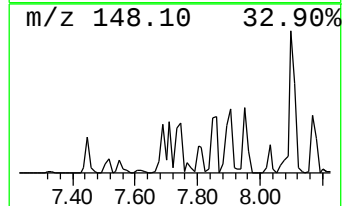
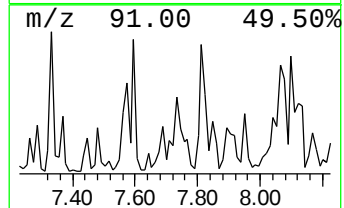
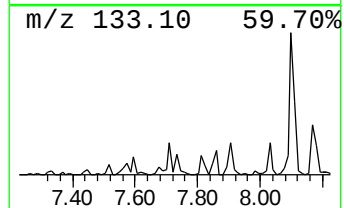
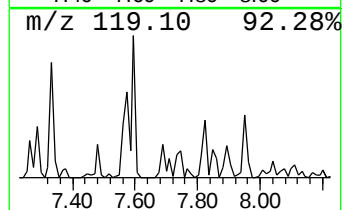
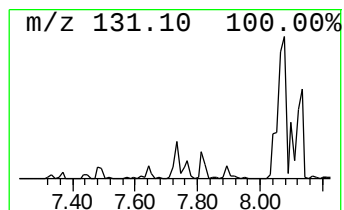
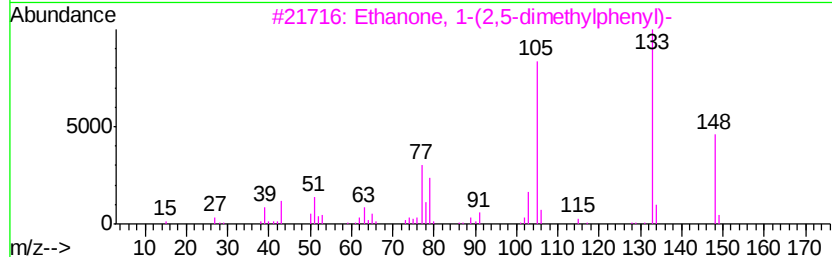
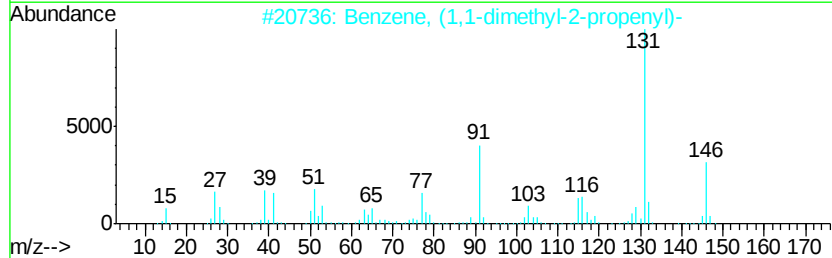
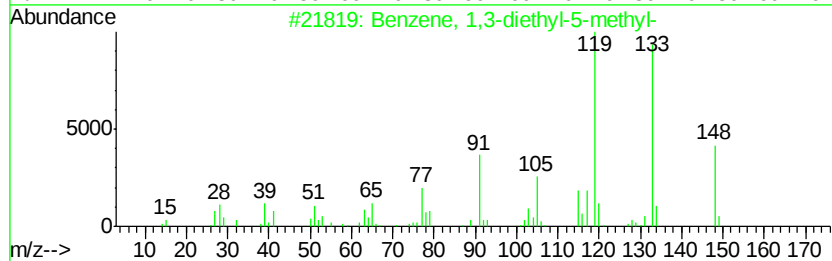
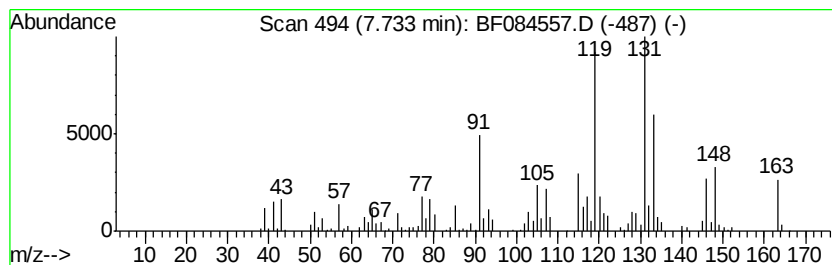
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	7.85 ng	1258860	Napthalene-d8	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	86
2		Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	83
3		Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	45
4		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	42
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
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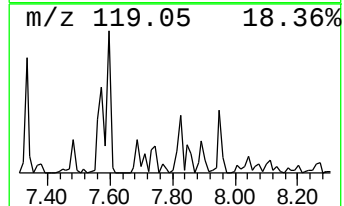
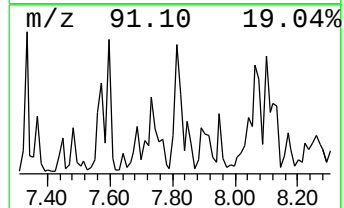
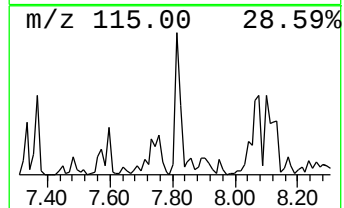
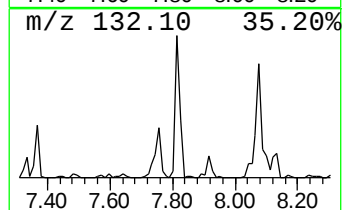
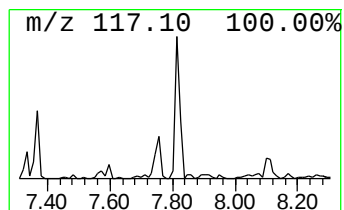
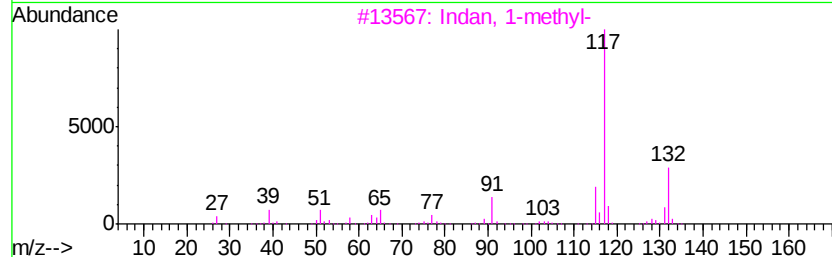
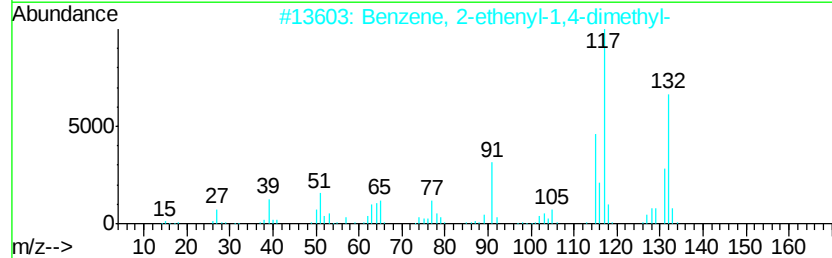
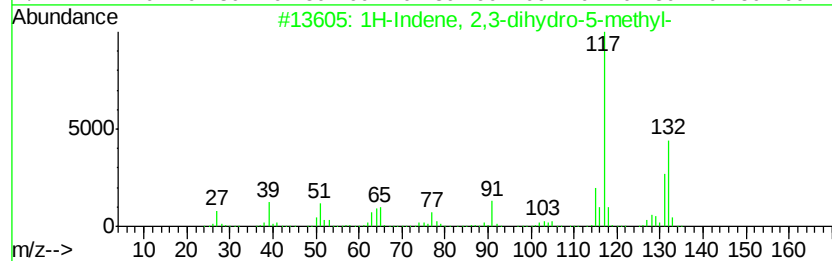
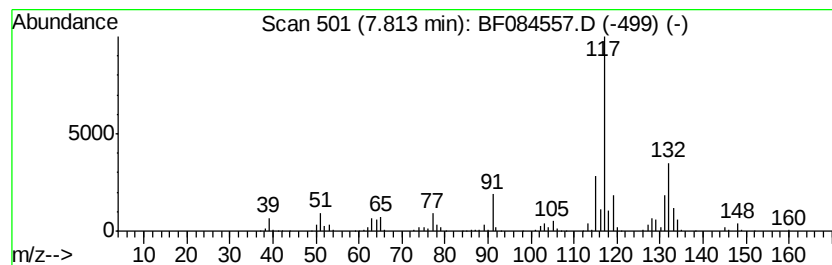
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	9.30 ng	1491490	Napthalene-d8	8.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3		Indan, 1-methyl-	132	C10H12	000767-58-8	87
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
Sample : H1169-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-004(0-2)

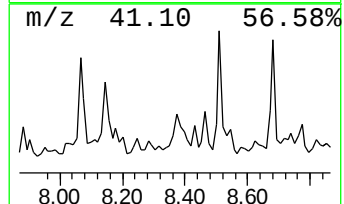
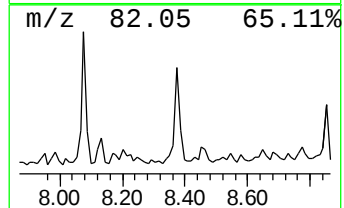
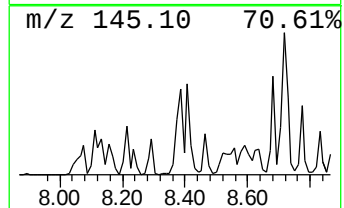
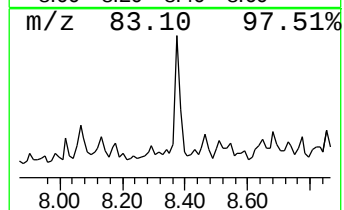
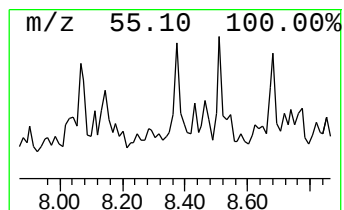
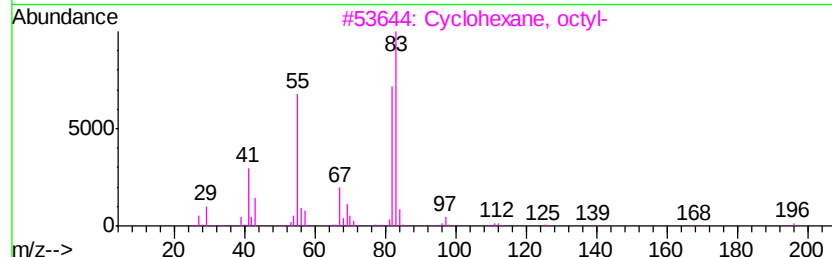
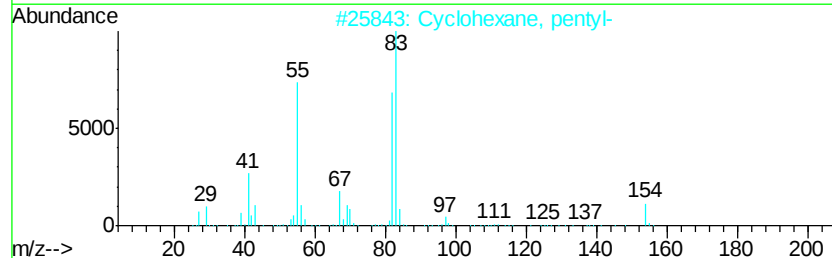
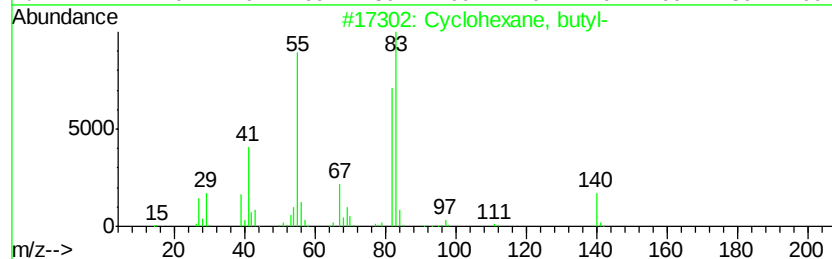
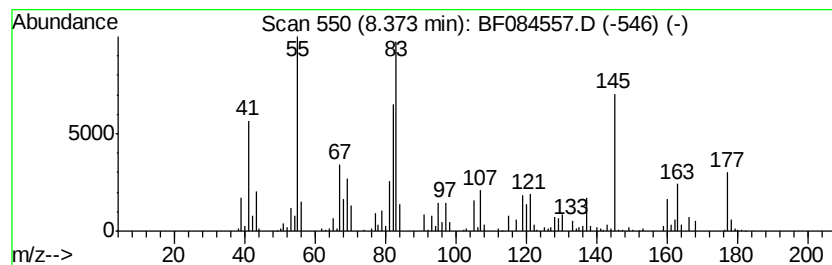
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown8.37 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.37	8.79 ng	1409670	Napthalene-d8	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	35
2			Cyclohexane, pentyl-	154	C11H22	004292-92-6	35
3			Cyclohexane, octyl-	196	C14H28	001795-15-9	35
4			Heptylcyclohexane	182	C13H26	005617-41-4	35
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
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Acq On : 19 Jan 2016 17:34
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Misc :
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Instrument :
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ClientSampled :
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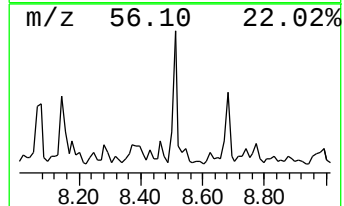
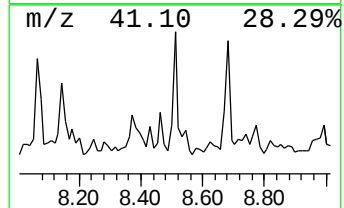
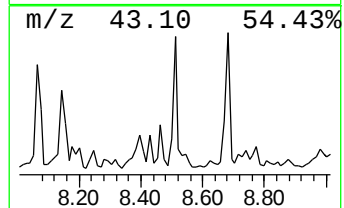
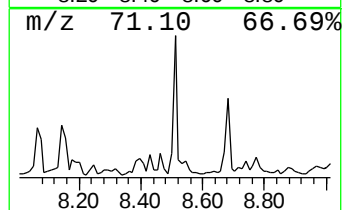
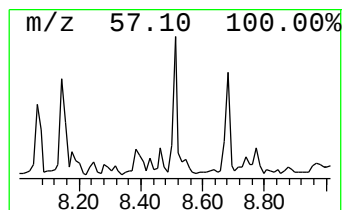
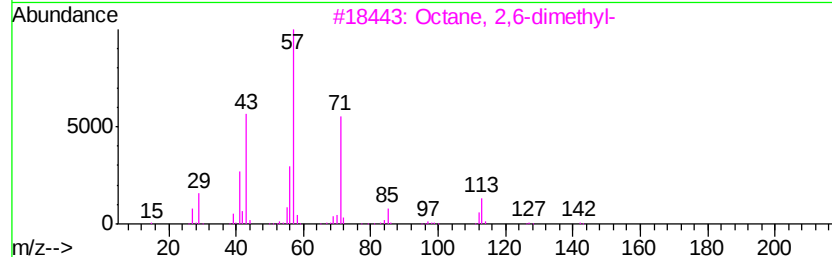
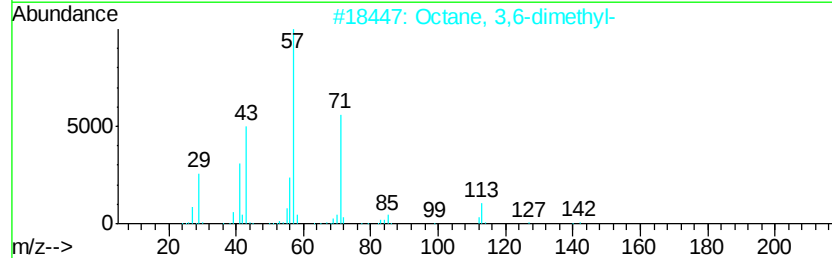
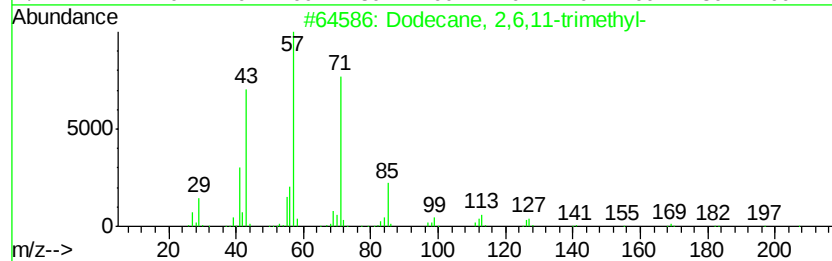
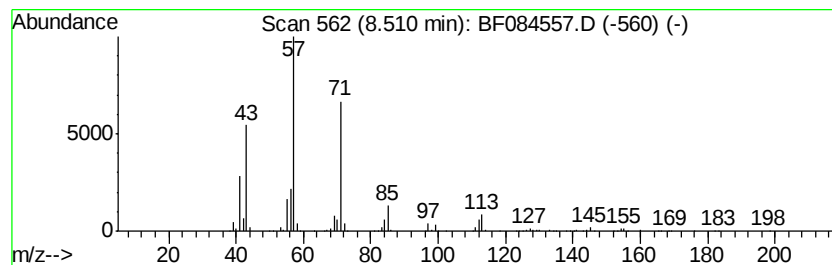
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dodecane, 2,6,11-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.51	7.49 ng	1201320	Napthalene-d8	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	78
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
3		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
4		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
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Sample : H1169-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

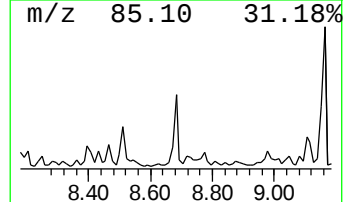
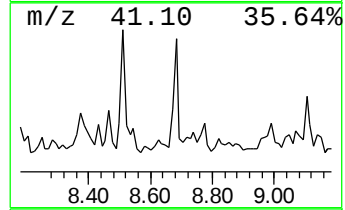
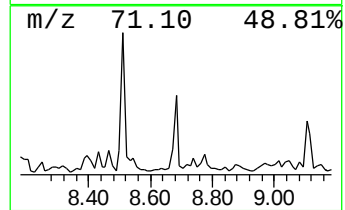
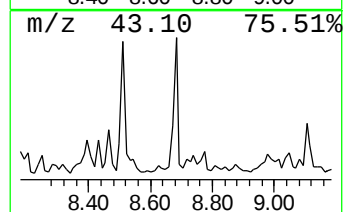
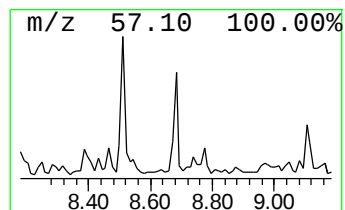
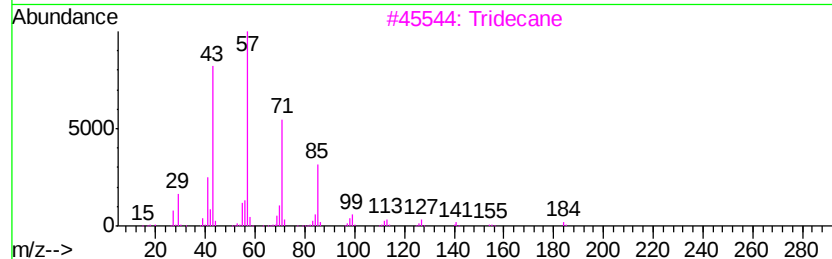
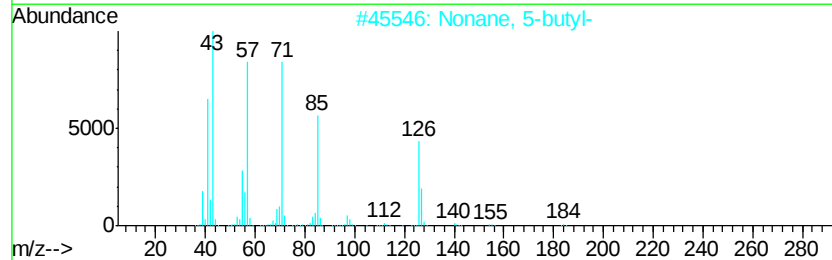
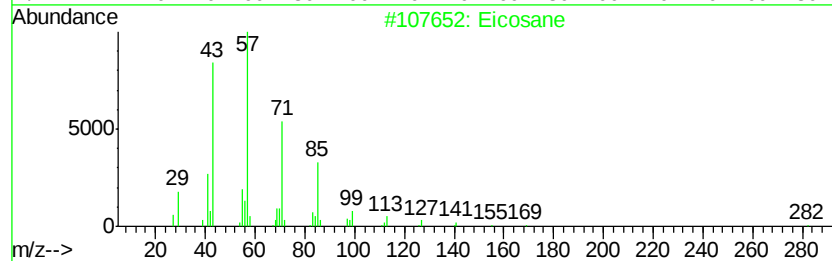
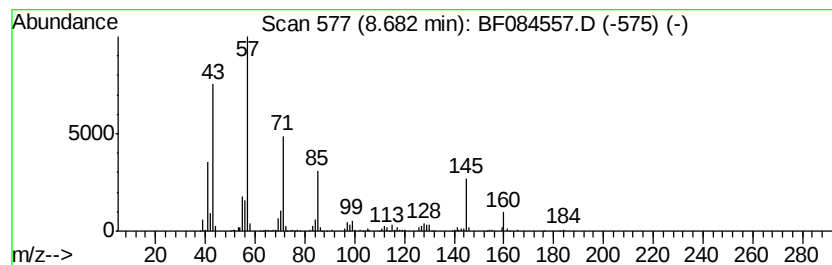
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.68	7.07 ng	1133620	Napthalene-d8	8.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	68
2	Nonane, 5-butyl-	184 C13H28	017312-63-9	60
3	Tridecane	184 C13H28	000629-50-5	58
4	Tetradecane	198 C14H30	000629-59-4	58
5	Decane, 2,3,5-trimethyl-	184 C13H28	062238-11-3	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
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Instrument :
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ClientSampleId :
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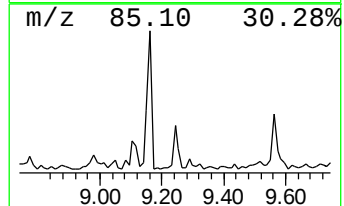
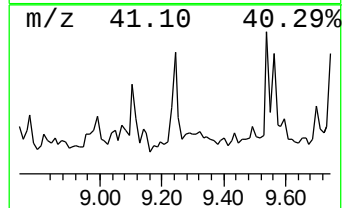
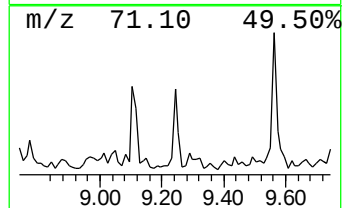
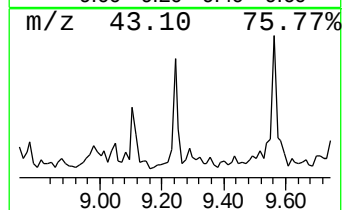
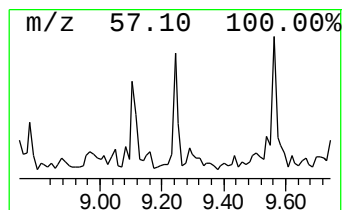
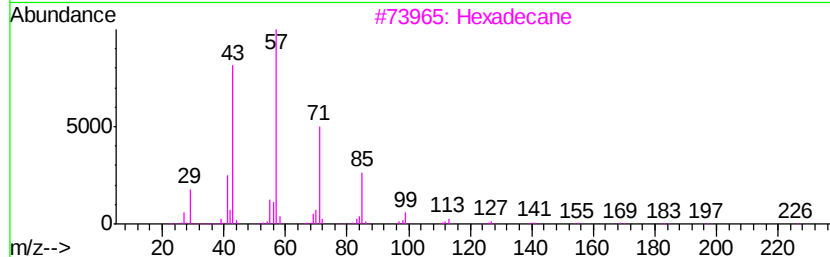
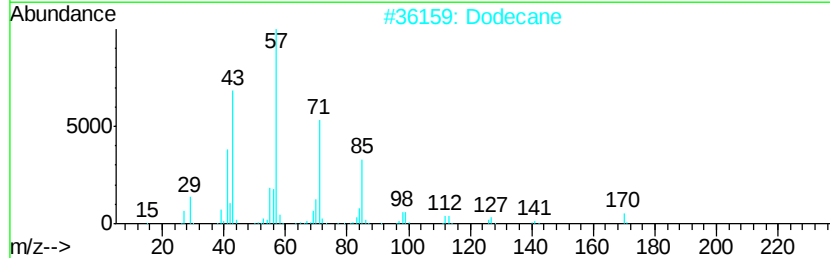
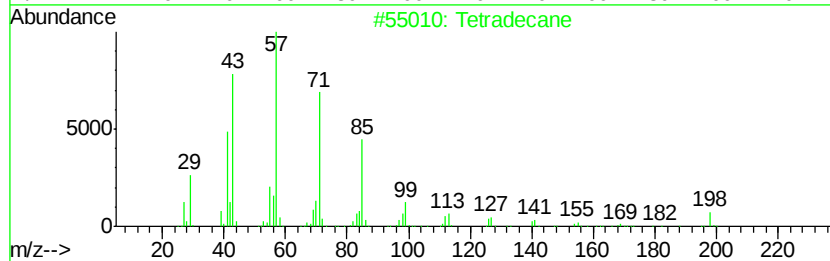
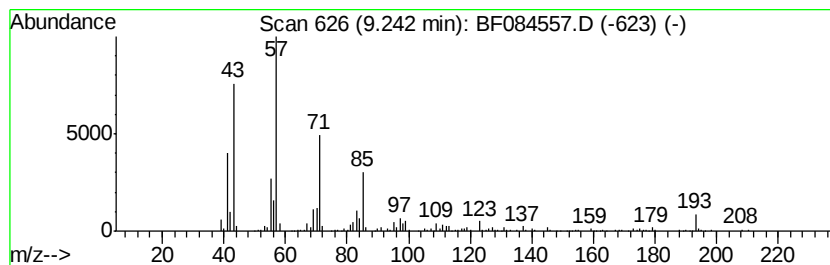
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetradecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.24	8.20 ng	910245	Acenaphthene-d10	9.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	93
2		Dodecane	170	C12H26	000112-40-3	72
3		Hexadecane	226	C16H34	000544-76-3	72
4		Eicosane	282	C20H42	000112-95-8	72
5		Pentadecane	212	C15H32	000629-62-9	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
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ALS Vial : 12 Sample Multiplier: 1

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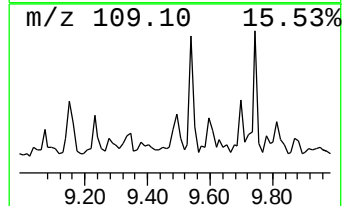
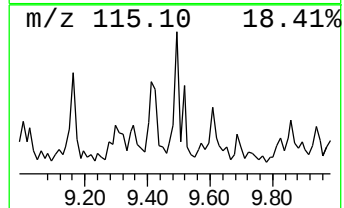
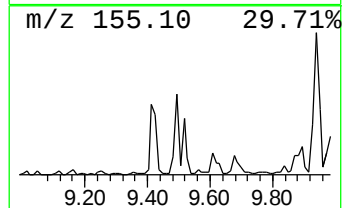
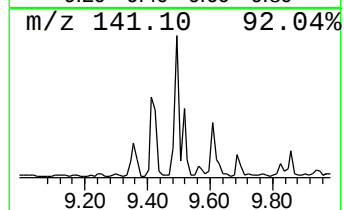
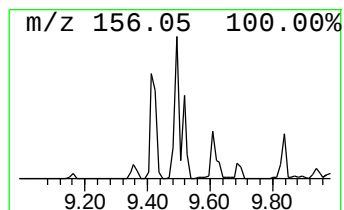
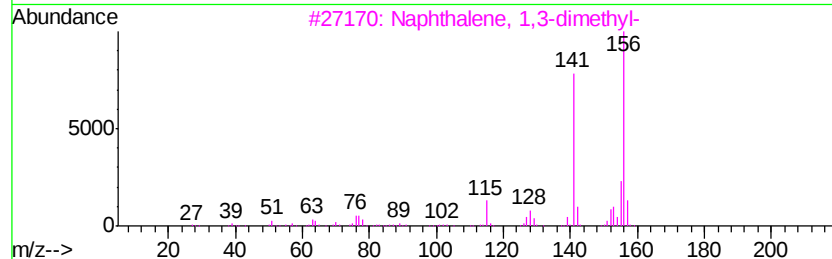
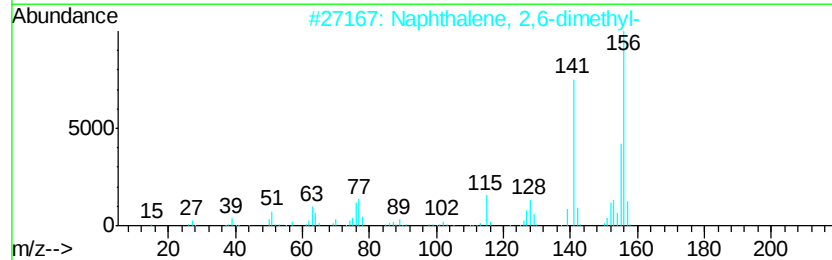
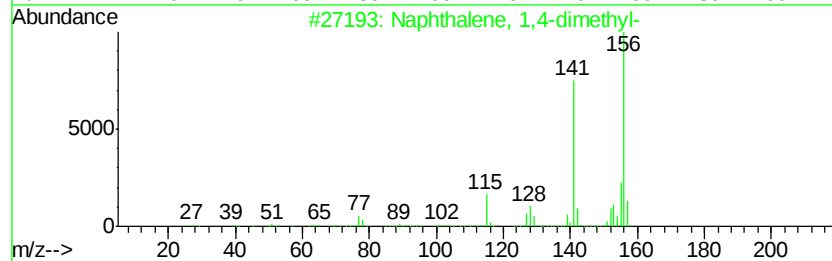
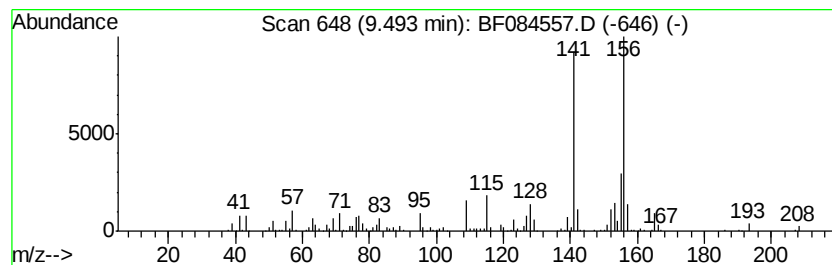
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 1,4-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	6.76 ng	750729	Acenaphthene-d10	9.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
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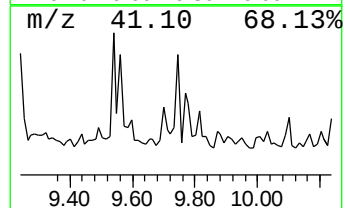
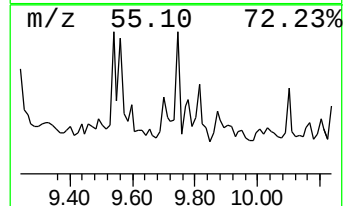
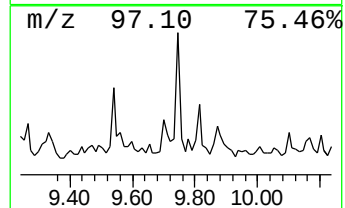
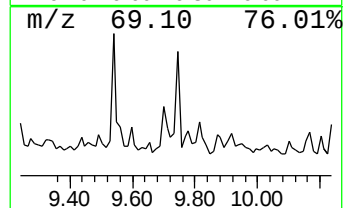
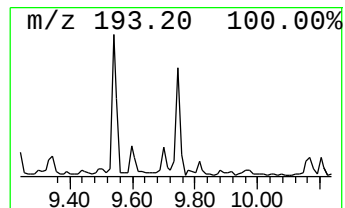
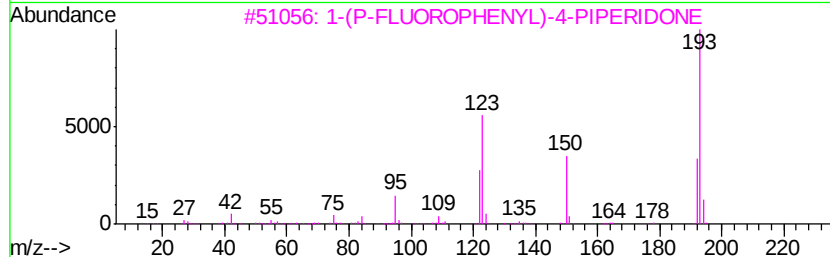
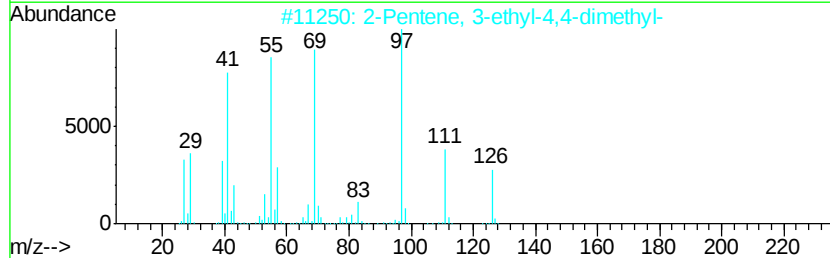
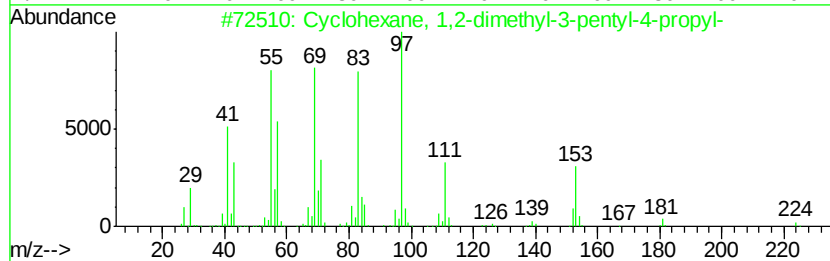
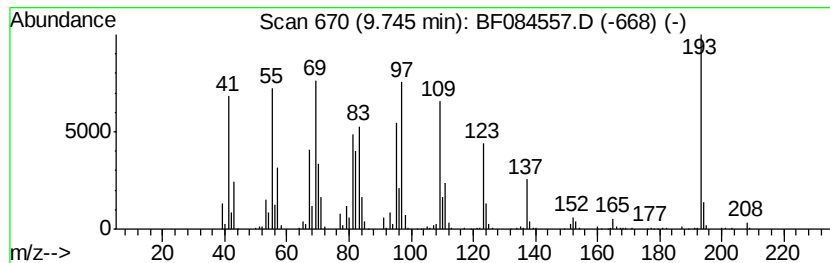
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown9.74 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.74	5.63 ng	624879	Acenaphthene-d10	9.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	27
2		2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	27
3		1-(P-FLUOROPHENYL)-4-PIPERIDONE	193	C11H12FN0	1000238-56-7	25
4		Cyclopentanecarboxylic acid, 4-t...	310	C20H38O2	1000280-58-9	22
5		1-Hexyne, 3-ethoxy-3,4-dimethyl-	154	C10H18O	054244-91-6	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
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Misc :
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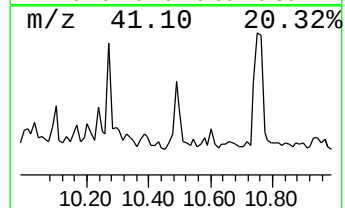
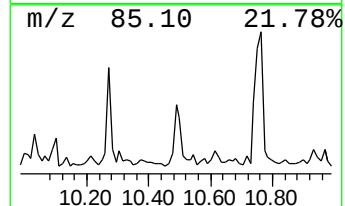
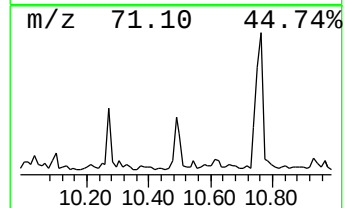
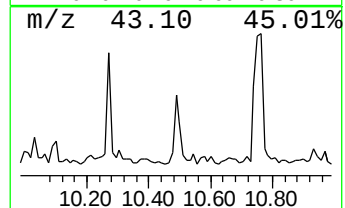
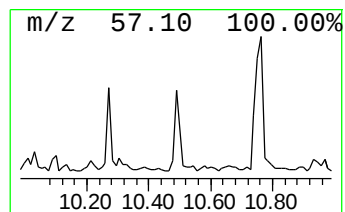
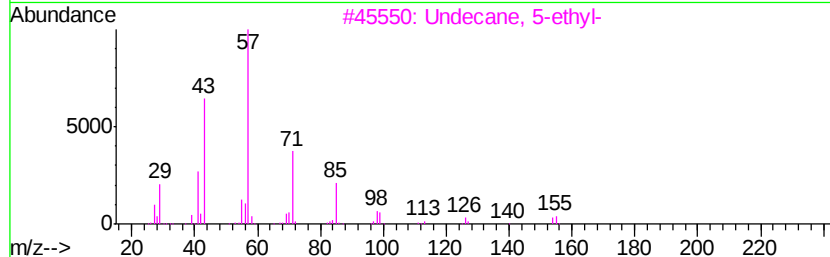
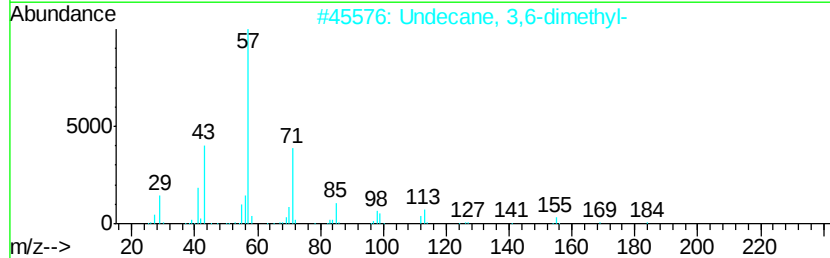
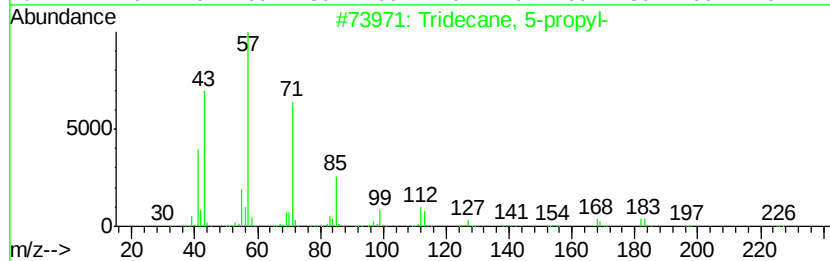
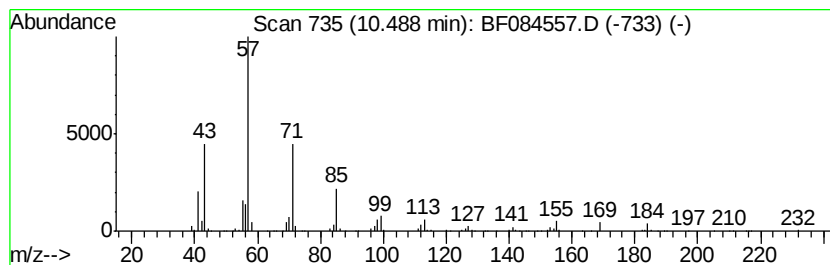
Instrument :
BNA_F
ClientSampleId :
SP-004(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Tridecane, 5-propyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.49	5.54 ng	614733	Acenaphthene-d10	9.84	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	80
2	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	72
3	Undecane, 5-ethyl-	184	C13H28	017453-94-0	72
4	Eicosane	282	C20H42	000112-95-8	64
5	Tridecane	184	C13H28	000629-50-5	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
Data File : BF084557.D
Acq On : 19 Jan 2016 17:34
Operator : SJ/UM
Sample : H1169-07
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-004(0-2)

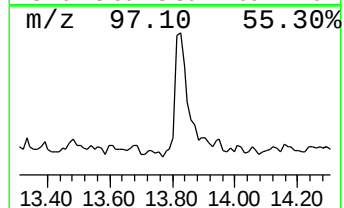
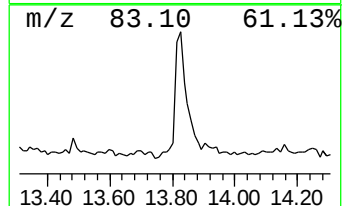
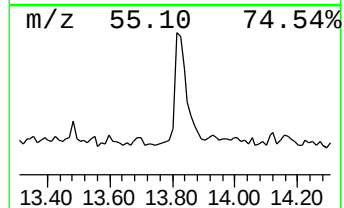
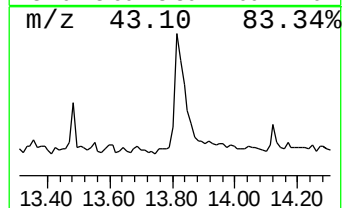
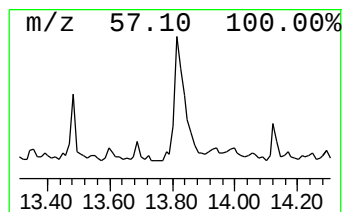
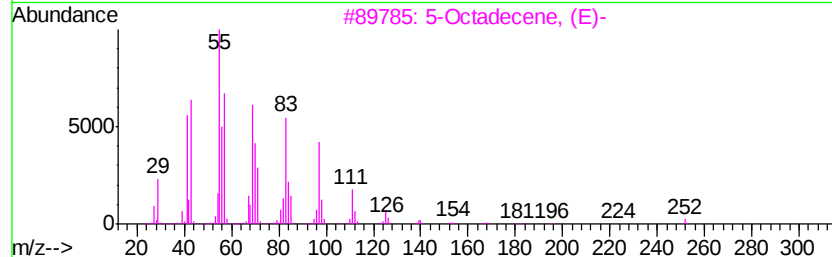
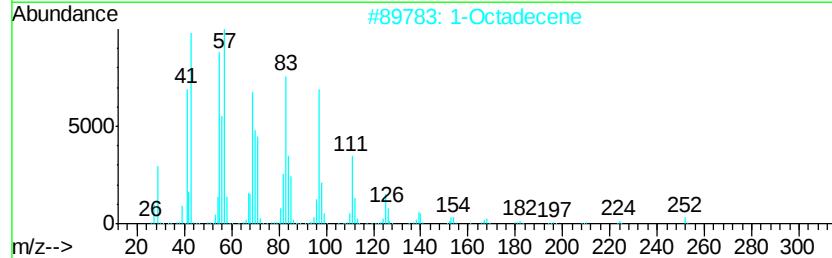
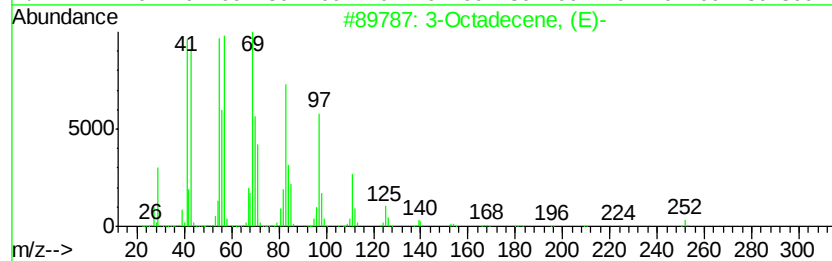
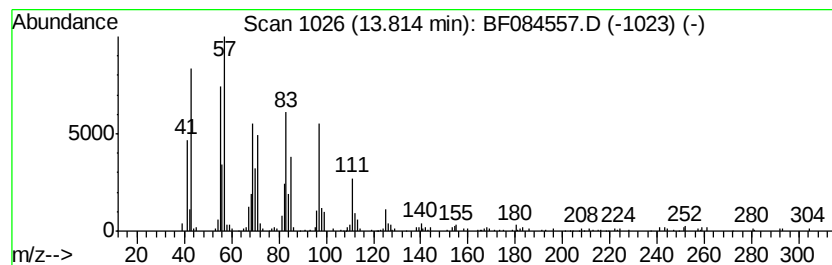
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 3-Octadecene, (E)- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	5.29 ng	416341	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Octadecene, (E)-	252	C18H36	007206-19-1	95
2			1-Octadecene	252	C18H36	000112-88-9	93
3			5-Octadecene, (E)-	252	C18H36	007206-21-5	92
4			5-Eicosene, (E)-	280	C20H40	074685-30-6	92
5			17-Pentatriacontene	491	C35H70	006971-40-0	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011916\
 Data File : BF084557.D
 Acq On : 19 Jan 2016 17:34
 Operator : SJ/UM
 Sample : H1169-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-004(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.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...	4.97	38.4	ng	3369000	1	6.80	1754490	20.0
Cyclohexane, 1-et...	5.80	4.9	ng	427947	1	6.80	1754490	20.0
unknown6.03	6.03	6.3	ng	554088	1	6.80	1754490	20.0
unknown6.57	6.57	65.3	ng	5726430	1	6.80	1754490	20.0
Decane	6.64	7.1	ng	623796	1	6.80	1754490	20.0
unknown7.07	7.07	7.2	ng	633320	1	6.80	1754490	20.0
Benzene, 2-ethyl-...	7.12	5.3	ng	464928	1	6.80	1754490	20.0
Benzene, 1,2,4,5-...	7.57	5.5	ng	890380	2	8.08	3206810	20.0
Benzene, 1,3-diet...	7.73	7.8	ng	1258860	2	8.08	3206810	20.0
1H-Indene, 2,3-di...	7.81	9.3	ng	1491490	2	8.08	3206810	20.0
unknown8.37	8.37	8.8	ng	1409670	2	8.08	3206810	20.0
Dodecane, 2,6,11-...	8.51	7.5	ng	1201320	2	8.08	3206810	20.0
Eicosane	8.68	7.1	ng	1133620	2	8.08	3206810	20.0
Tetradecane	9.24	8.2	ng	910245	3	9.84	2219710	20.0
Naphthalene, 1,4-...	9.49	6.8	ng	750729	3	9.84	2219710	20.0
unknown9.74	9.74	5.6	ng	624879	3	9.84	2219710	20.0
Tridecane, 5-propyl-	10.49	5.5	ng	614733	3	9.84	2219710	20.0
3-Octadecene, (E)-	13.81	5.3	ng	416341	5	13.95	1575540	20.0