

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
 Data File : BF088212.D
 Acq On : 21 Jun 2016 11:23
 Operator : UM/SJ
 Sample : H3580-21
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 08-B6(8-16)C

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-BF060716.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.655	5	6	14	rVB	124430	196765	9.41%	0.833%
2	1.770	14	16	53	rVB	1039265	2090143	100.00%	8.852%
3	4.821	281	283	287	rBV	62648	100994	4.83%	0.428%
4	5.267	320	322	331	rBV	1482932	1590281	76.08%	6.735%
5	5.656	350	356	358	rBV2	23363	41365	1.98%	0.175%
6	5.793	366	368	371	rVB	24619	32316	1.55%	0.137%
7	5.861	371	374	376	rBV2	25725	47732	2.28%	0.202%
8	5.907	376	378	381	rVV2	86703	134103	6.42%	0.568%
9	5.964	381	383	384	rVV	48671	67829	3.25%	0.287%
10	5.987	384	385	391	rVB4	26420	69502	3.33%	0.294%
11	6.101	391	395	398	rBV	46162	87891	4.21%	0.372%
12	6.181	400	402	403	rBV	52805	49402	2.36%	0.209%
13	6.216	403	405	407	rVB2	32053	44939	2.15%	0.190%
14	6.273	407	410	412	rBV2	25116	54019	2.58%	0.229%
15	6.330	412	415	417	rVV	1615033	1792278	85.75%	7.590%
16	6.410	419	422	423	rVV	55222	110861	5.30%	0.469%
17	6.444	423	425	427	rVV	1746084	1799443	86.09%	7.621%
18	6.524	429	432	433	rVB2	42406	68404	3.27%	0.290%
19	6.604	433	439	440	rBV3	44321	80661	3.86%	0.342%
20	6.639	440	442	443	rVV2	41486	64067	3.07%	0.271%
21	6.673	443	445	448	rVB	341249	498074	23.83%	2.109%
22	6.730	448	450	451	rBV	48523	64108	3.07%	0.271%
23	6.764	451	453	454	rVV2	48278	72634	3.48%	0.308%
24	6.787	454	455	456	rVV	72424	95253	4.56%	0.403%
25	6.822	456	458	462	rVB	1482067	1426713	68.26%	6.042%
26	6.902	462	465	467	rBV3	43521	119075	5.70%	0.504%
27	6.936	467	468	471	rVB2	60663	81769	3.91%	0.346%
28	7.016	474	475	477	rBV2	33101	59290	2.84%	0.251%
29	7.062	477	479	480	rVV	53123	56762	2.72%	0.240%
30	7.084	480	481	486	rVV5	33577	82839	3.96%	0.351%
31	7.210	488	492	493	rVV2	67343	148433	7.10%	0.629%
32	7.233	493	494	497	rVV	813884	938396	44.90%	3.974%
33	7.313	499	501	504	rBV3	56573	108616	5.20%	0.460%
34	7.393	504	508	510	rVB4	23179	62201	2.98%	0.263%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	7.439	510	512	514	rBV3	14004	33362	1.60%	0.141%
36	7.473	514	515	517	rVB	94078	81178	3.88%	0.344%
37	7.519	517	519	521	rBV3	23643	37547	1.80%	0.159%
38	7.587	521	525	528	rVB4	73300	152921	7.32%	0.648%
39	7.667	530	532	534	rVB3	19403	30827	1.47%	0.131%
40	7.713	534	536	540	rBV2	32012	69420	3.32%	0.294%
41	7.782	540	542	544	rVB	41615	41273	1.97%	0.175%
42	7.850	544	548	550	rBV4	25470	80741	3.86%	0.342%
43	7.907	550	553	554	rBV3	34252	59710	2.86%	0.253%
44	7.953	555	557	559	rVV	751400	638385	30.54%	2.704%
45	8.022	562	563	565	rVB	22361	26600	1.27%	0.113%
46	8.124	570	572	573	rBV	24057	29064	1.39%	0.123%
47	8.159	573	575	577	rBV	41725	40766	1.95%	0.173%
48	8.250	579	583	584	rBV2	55261	85348	4.08%	0.361%
49	8.387	593	595	599	rVB	279109	316793	15.16%	1.342%
50	8.456	599	601	603	rVB2	27165	27486	1.32%	0.116%
51	8.490	603	604	608	rBV4	34182	72346	3.46%	0.306%
52	8.547	608	609	612	rBV2	34492	66562	3.18%	0.282%
53	8.650	616	618	620	rVB	60449	74713	3.57%	0.316%
54	8.696	620	622	623	rBV	26254	29150	1.39%	0.123%
55	8.753	626	627	631	rVB3	31783	52060	2.49%	0.220%
56	8.833	631	634	635	rBV2	47545	61796	2.96%	0.262%
57	8.867	635	637	639	rVB2	64688	71546	3.42%	0.303%
58	8.959	644	645	646	rVV	47108	55468	2.65%	0.235%
59	8.982	646	647	649	rVV	243523	217355	10.40%	0.920%
60	9.027	649	651	653	rVV	1587242	1549960	74.16%	6.564%
61	9.130	655	660	661	rVV2	81461	187084	8.95%	0.792%
62	9.165	661	663	665	rVV2	48413	105431	5.04%	0.446%
63	9.233	667	669	670	rVV2	34460	36942	1.77%	0.156%
64	9.279	670	673	674	rVV	20654	38999	1.87%	0.165%
65	9.302	674	675	676	rVV	40645	39221	1.88%	0.166%
66	9.370	679	681	682	rVV	51678	79733	3.81%	0.338%
67	9.405	682	684	685	rVV2	83411	142713	6.83%	0.604%
68	9.427	685	686	688	rVV2	672181	698624	33.42%	2.959%
69	9.496	690	692	693	rVB2	35630	35027	1.68%	0.148%
70	9.610	701	702	704	rVB	73790	55406	2.65%	0.235%
71	9.645	704	705	707	rVV	75465	62546	2.99%	0.265%

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72	9.702	707	710	712 rVV	714599	689993	33.01%	2.922%
73	9.748	712	714	716 rVV2	22917	33371	1.60%	0.141%
74	9.816	718	720	722 rVB	38330	46051	2.20%	0.195%
75	9.908	722	728	729 rBV3	57780	148299	7.10%	0.628%
76	9.965	732	733	735 rVB	85631	70962	3.40%	0.301%
77	10.102	743	745	747 rBV3	21030	33797	1.62%	0.143%
78	10.136	747	748	750 rVB	46988	59158	2.83%	0.251%
79	10.365	765	768	770 rBV	115868	161608	7.73%	0.684%
80	10.490	776	779	781 rBV	667306	751087	35.93%	3.181%
81	10.570	784	786	788 rVB2	26429	35058	1.68%	0.148%
82	10.628	788	791	793 rBV	217902	318074	15.22%	1.347%
83	10.810	805	807	811 rVB4	25337	65840	3.15%	0.279%
84	10.936	816	818	820 rBV2	24549	41589	1.99%	0.176%
85	11.062	827	829	830 rBV	40162	55461	2.65%	0.235%
86	11.085	830	831	834 rVB	96811	91282	4.37%	0.387%
87	11.176	837	839	843 rBV	580193	518814	24.82%	2.197%
88	11.771	890	891	897 rVB5	25661	65533	3.14%	0.278%
89	12.113	919	921	923 rVB2	26273	35622	1.70%	0.151%
90	12.171	923	926	928 rBV4	21354	47038	2.25%	0.199%
91	12.228	929	931	933 rBV2	22673	28663	1.37%	0.121%
92	12.388	943	945	946 rBV2	42181	46206	2.21%	0.196%
93	12.422	946	948	951 rVV4	29234	36416	1.74%	0.154%
94	12.616	963	965	966 rBV	33301	36379	1.74%	0.154%
95	12.754	975	977	980 rBV	886778	1022383	48.91%	4.330%
96	12.891	987	989	991 rBV	40015	37911	1.81%	0.161%
97	13.691	1057	1059	1067 rVB4	49935	108268	5.18%	0.459%
98	13.805	1067	1069	1073 rBV	513639	530605	25.39%	2.247%
99	15.177	1187	1189	1197 rBV	347000	546213	26.13%	2.313%
100	16.468	1300	1302	1308 rBV5	48162	129856	6.21%	0.550%

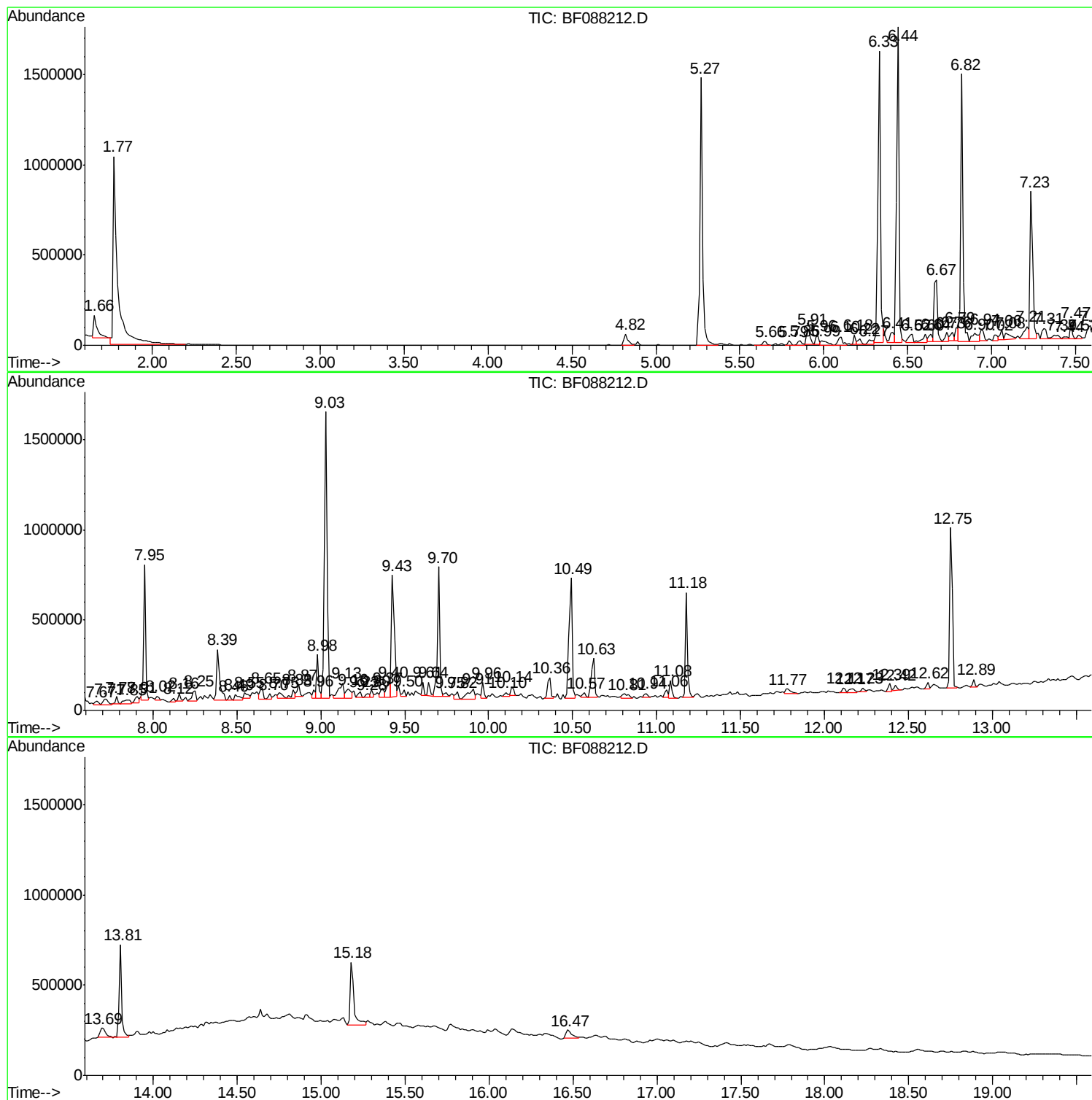
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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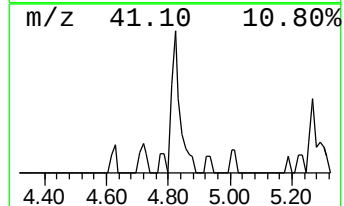
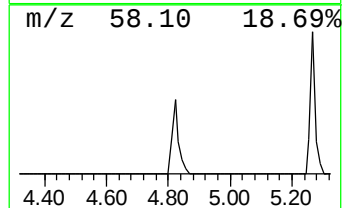
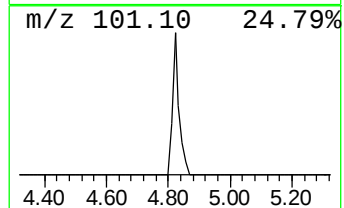
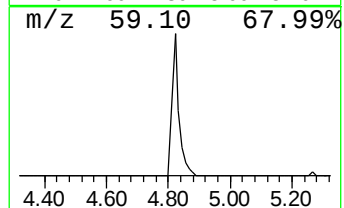
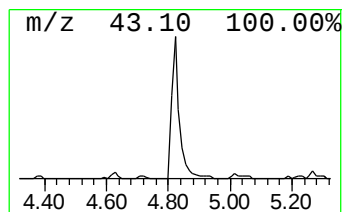
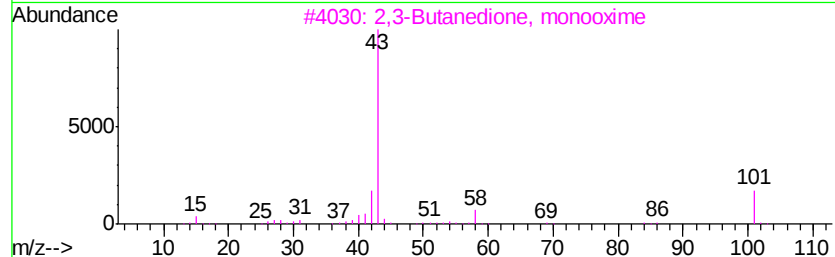
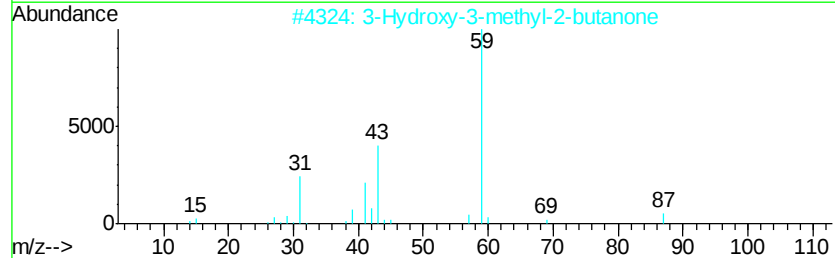
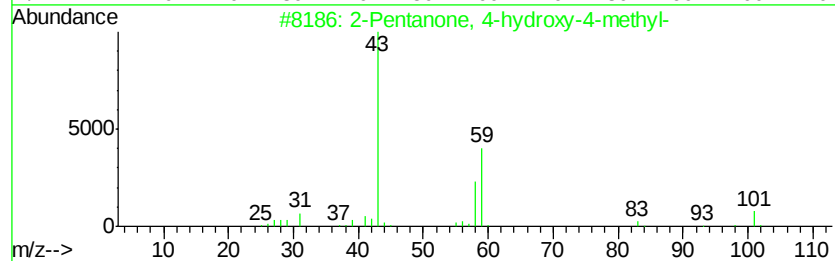
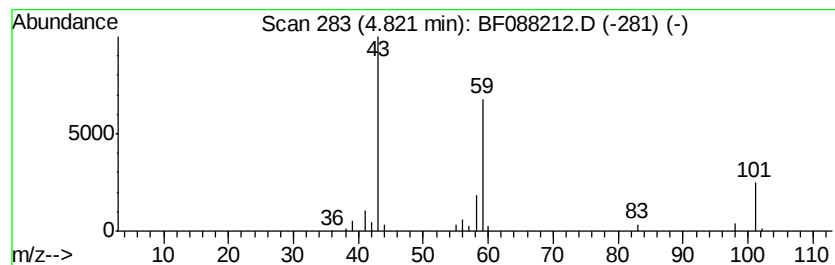
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	4.06 ng	100994	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4		5-Hexen-2-one	98	C6H10O	000109-49-9	10
5		Guanidine	59	CH5N3	000113-00-8	9



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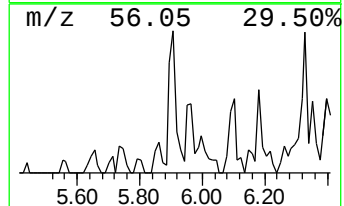
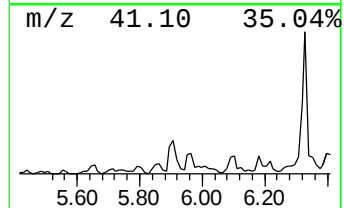
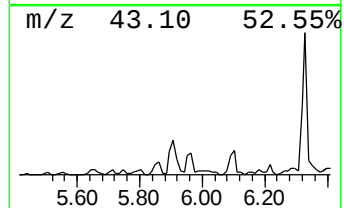
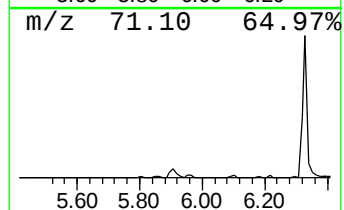
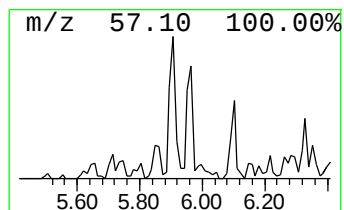
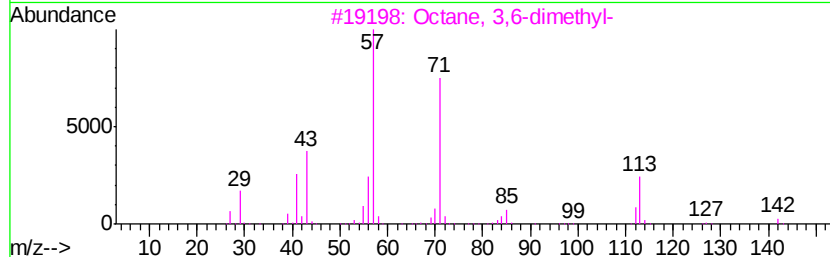
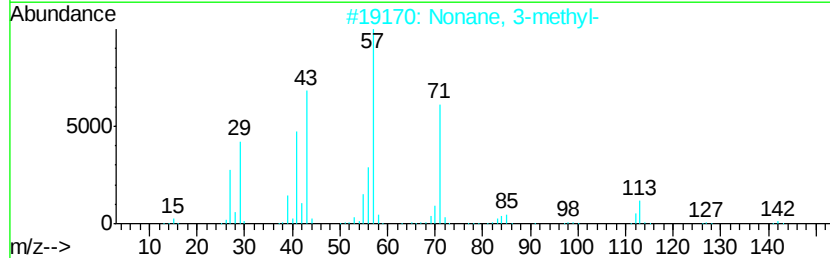
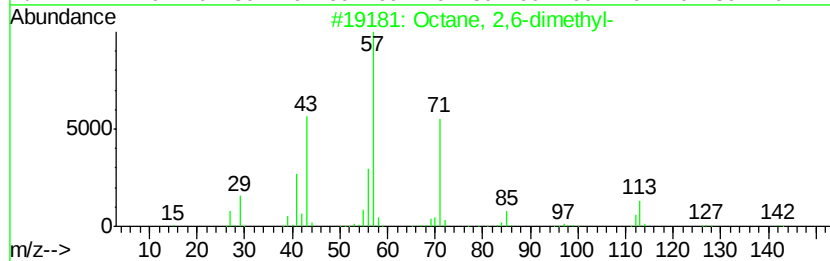
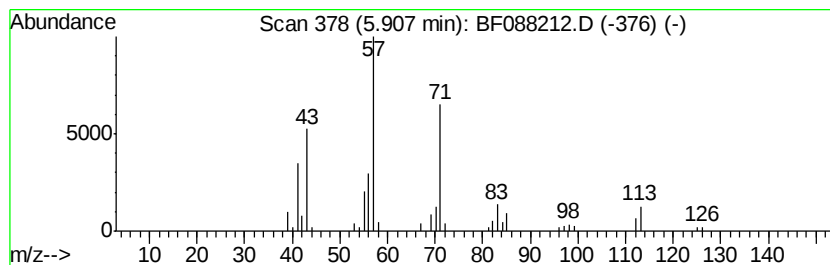
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.91	5.38 ng	134103	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	72
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
4		Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	64
5		Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	64



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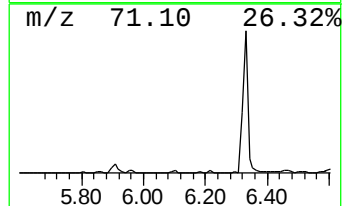
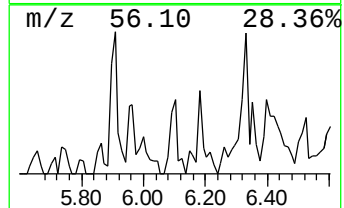
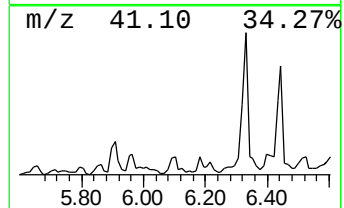
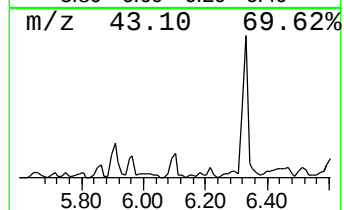
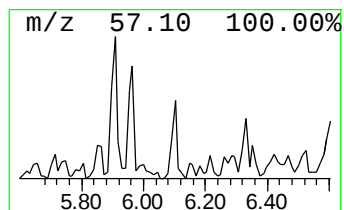
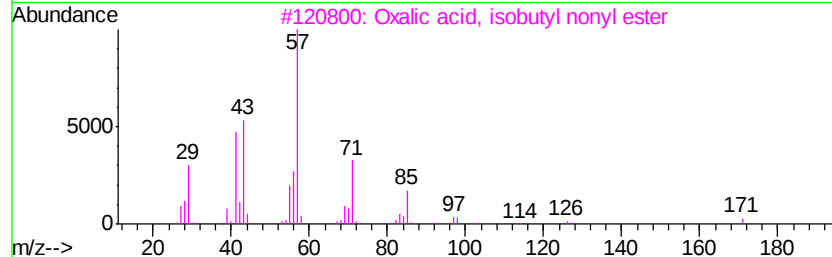
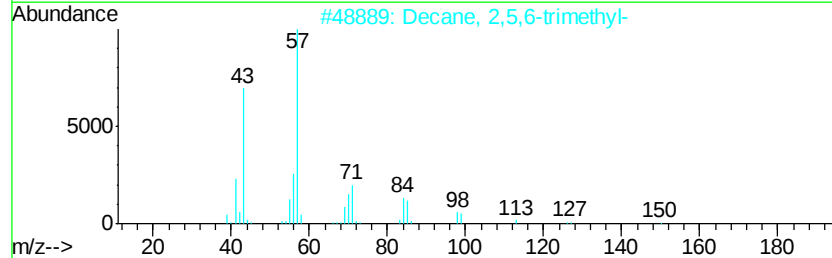
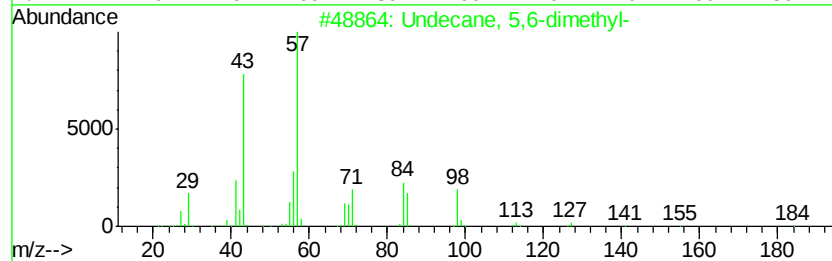
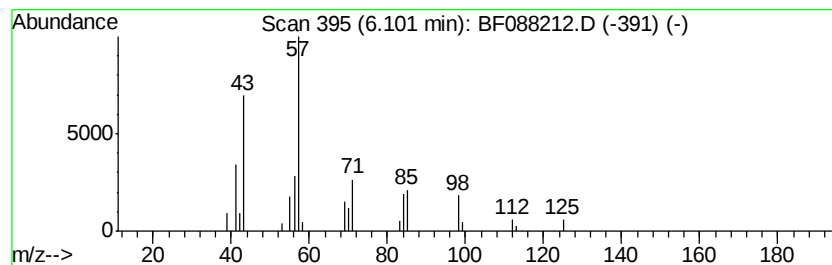
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Undecane, 5,6-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.10	3.53 ng	87891	1,4-Dichlorobenzene-d4	6.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 5,6-dimethyl-	184	C13H28	017615-91-7	72
2			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	50
3			Oxalic acid, isobutyl nonyl ester	272	C15H28O4	1000309-37-4	50
4			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	43
5			Undecane	156	C11H24	001120-21-4	43



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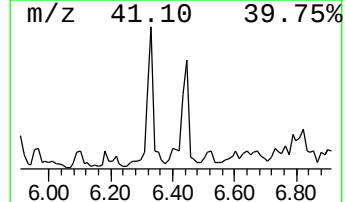
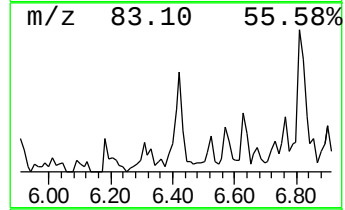
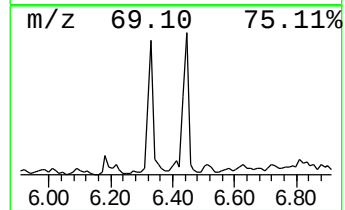
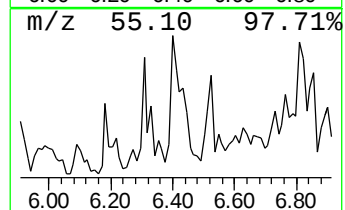
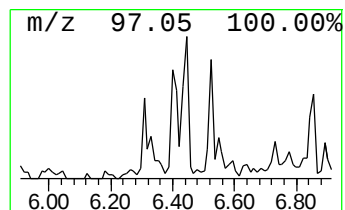
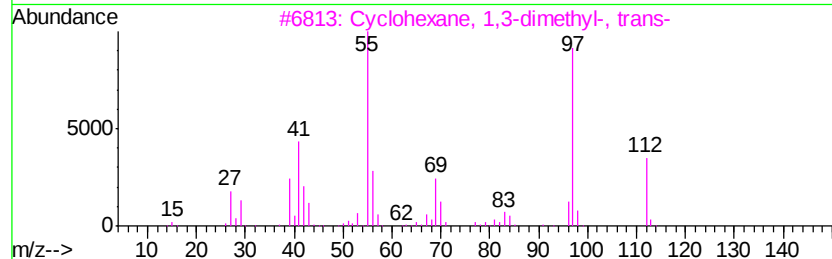
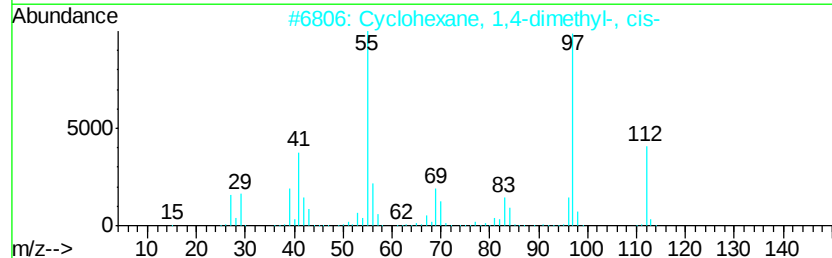
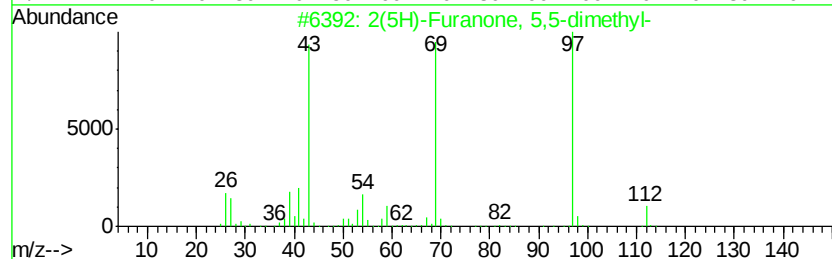
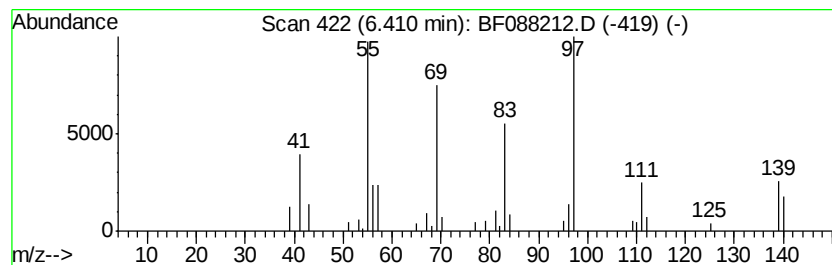
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown6.41 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	4.45 ng	110861	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(5H)-Furanone, 5,5-dimethyl-	112	C6H8O2	020019-64-1	38
2		Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	38
3		Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	38
4		Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	38
5		Sulfurous acid, cyclohexylmethyl...	262	C13H26O3S	1000309-21-5	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
Sample : H3580-21
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
08-B6(8-16)C

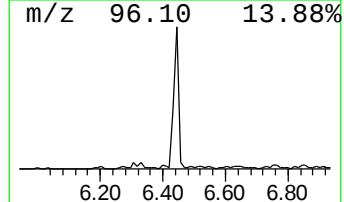
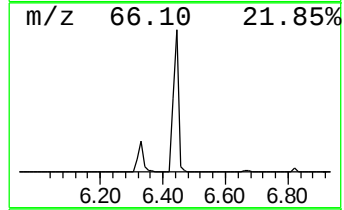
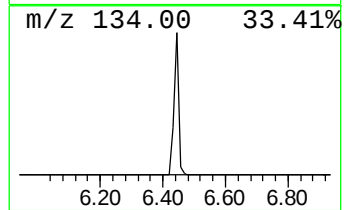
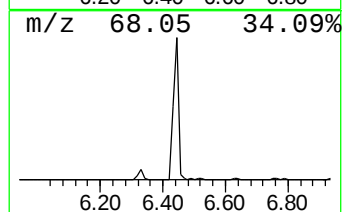
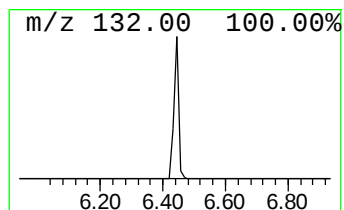
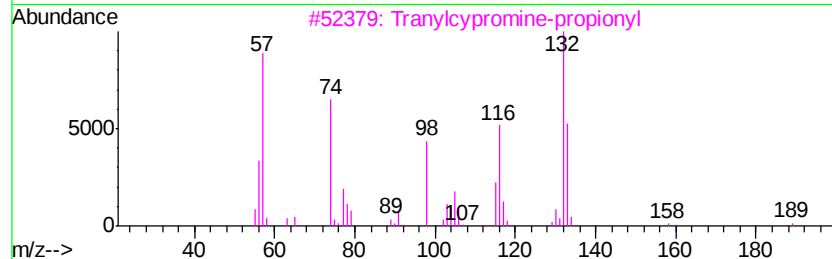
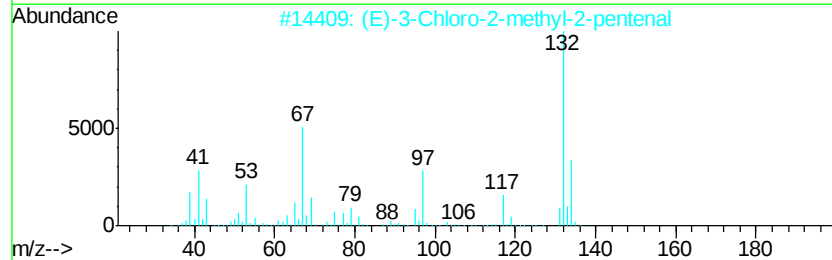
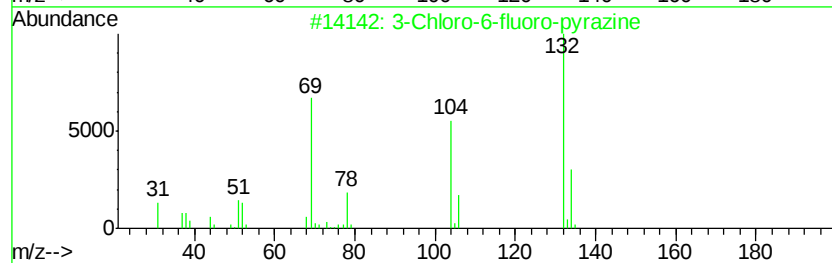
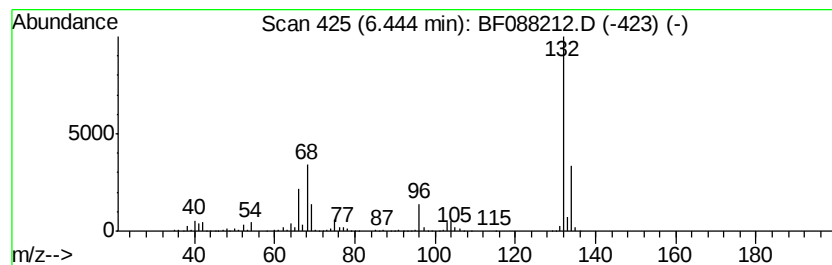
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.44 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.44	72.26 ng	1799440	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		Xenon	132	Xe	007440-63-3	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
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Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
08-B6(8-16)C

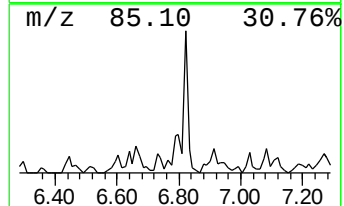
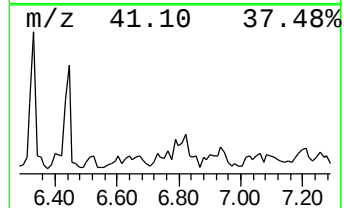
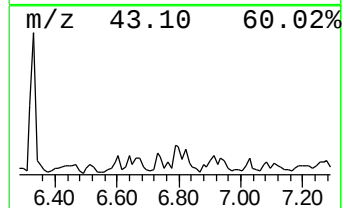
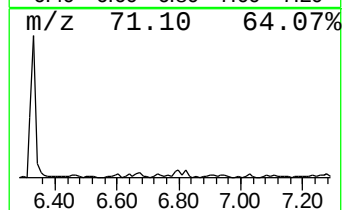
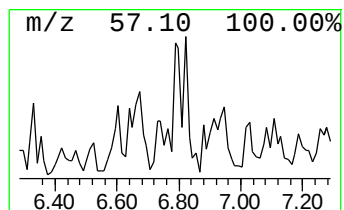
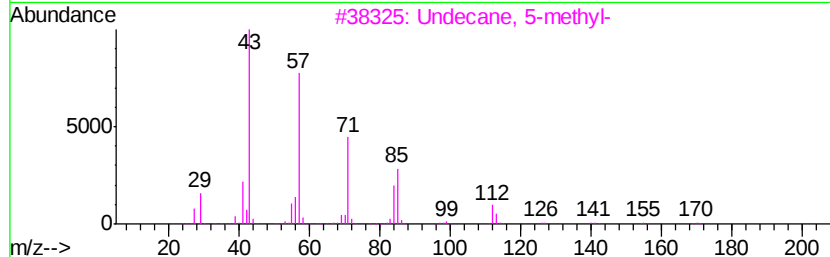
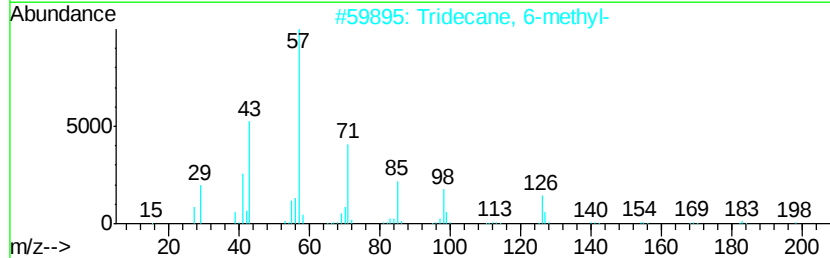
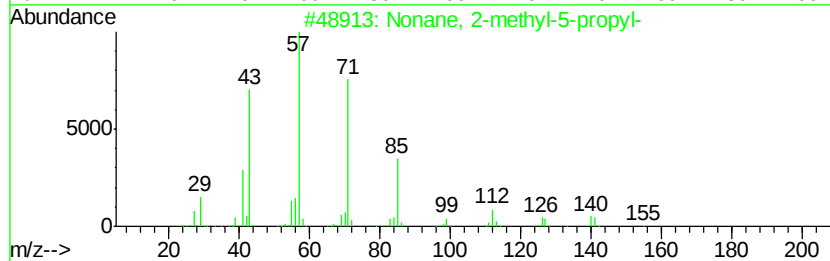
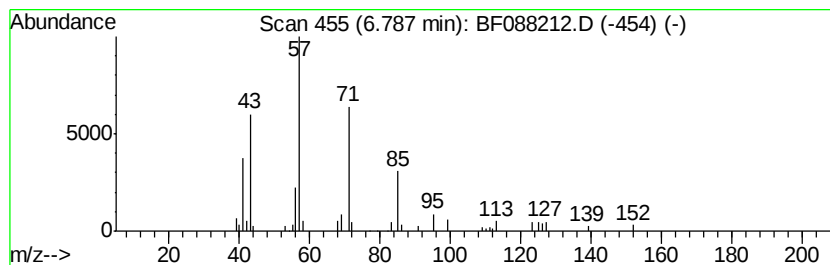
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonane, 2-methyl-5-propyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.79	3.82 ng	95253	1,4-Dichlorobenzene-d4	6.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	64
2		Tridecane, 6-methyl-	198	C14H30	013287-21-3	59
3		Undecane, 5-methyl-	170	C12H26	001632-70-8	59
4		Pentadecane	212	C15H32	000629-62-9	59
5		Hexadecane	226	C16H34	000544-76-3	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
Sample : H3580-21
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
08-B6(8-16)C

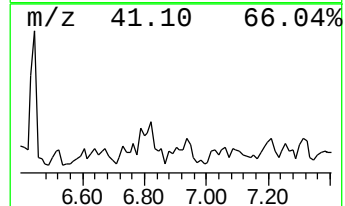
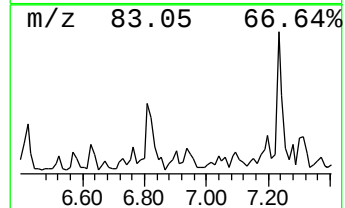
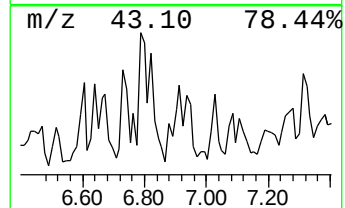
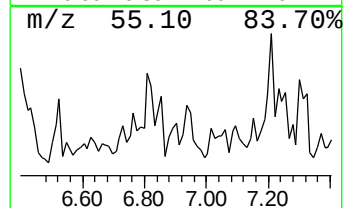
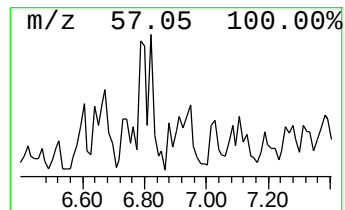
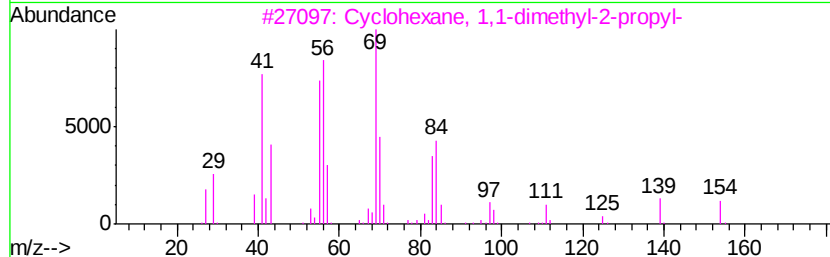
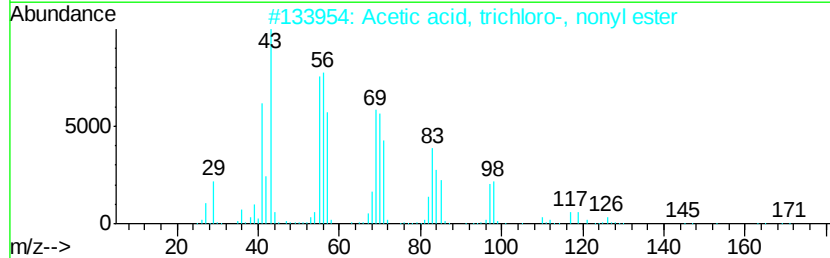
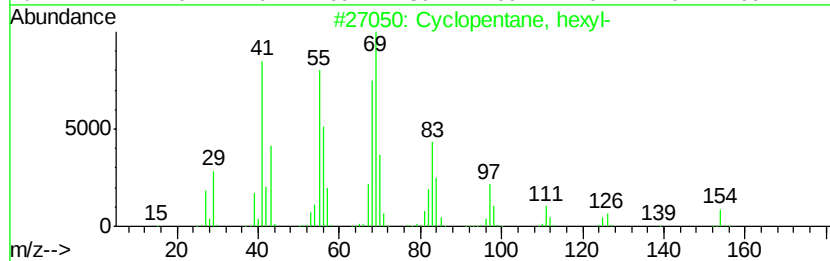
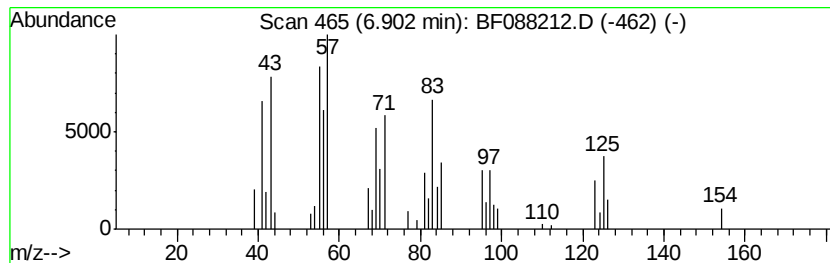
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown6.90 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.90	4.78 ng	119075	1,4-Dichlorobenzene-d4	6.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, hexyl-	154	C11H22	004457-00-5	38
2			Acetic acid, trichloro-, nonyl e...	288	C11H19Cl3O2	065611-32-7	35
3			Cyclohexane, 1,1-dimethyl-2-propyl-	154	C11H22	081983-71-3	25
4			(2-Methylbutyl)cyclohexane	154	C11H22	054105-77-0	25
5			Ethanone, 1-(2,2-dimethylcyclope...	140	C9H16O	003664-75-3	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
Sample : H3580-21
Misc :
ALS Vial : 40 Sample Multiplier: 1

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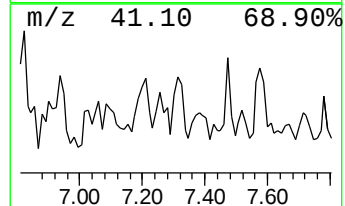
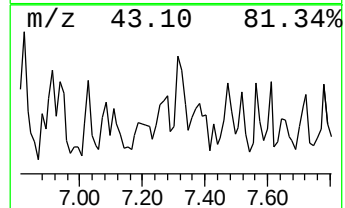
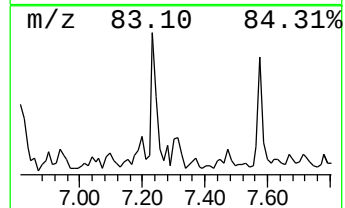
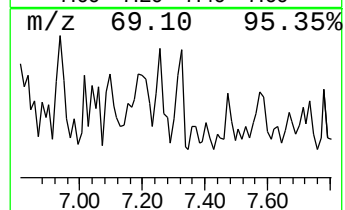
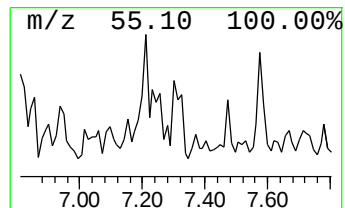
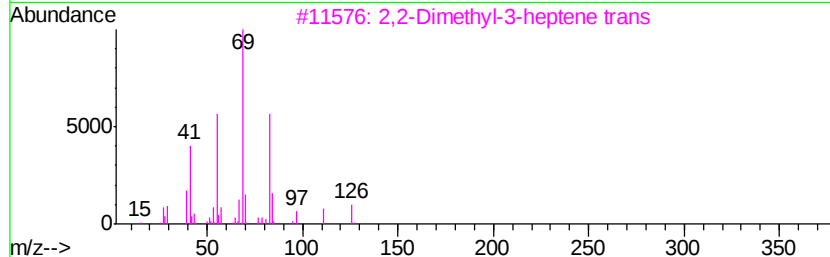
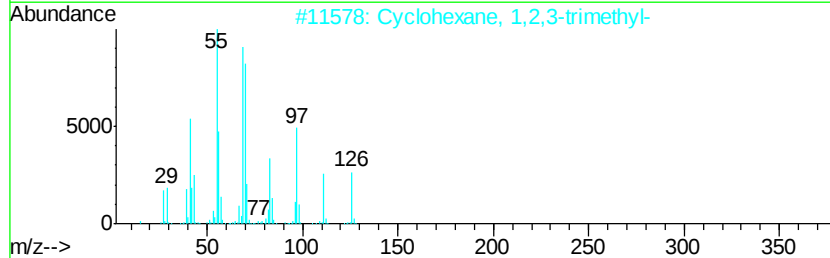
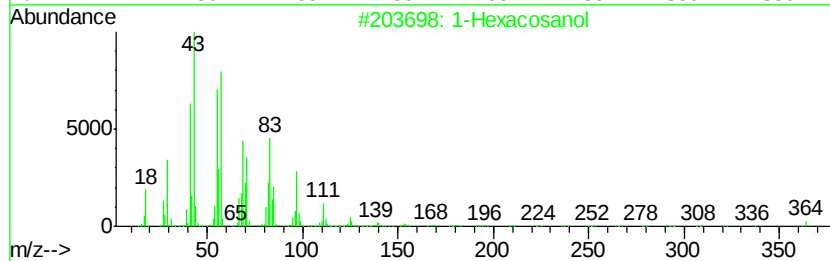
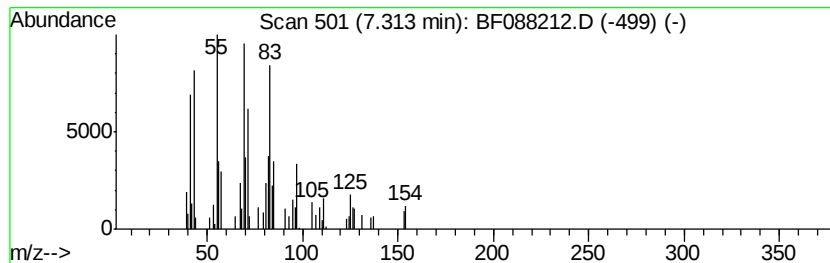
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown7.31 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.31	4.36 ng	108616	1,4-Dichlorobenzene-d4	6.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexacosanol	382	C26H54O	000506-52-5	43
2			Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	43
3			2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	38
4			Cyclohexane, 1,2-diethyl-3-methyl-	154	C11H22	061141-80-8	30
5			(2-Methylbutyl)cyclohexane	154	C11H22	054105-77-0	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
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Operator : UM/SJ
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Misc :
ALS Vial : 40 Sample Multiplier: 1

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ClientSampleId :
08-B6(8-16)C

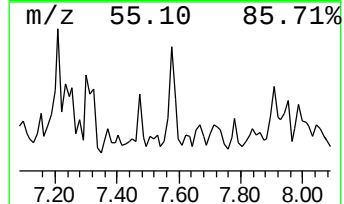
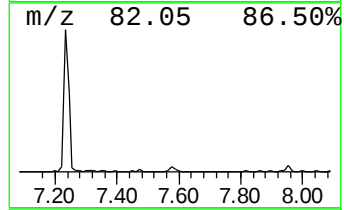
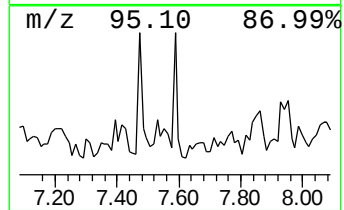
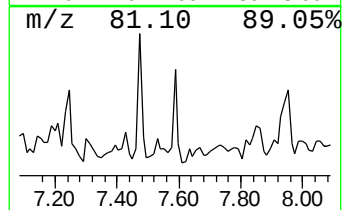
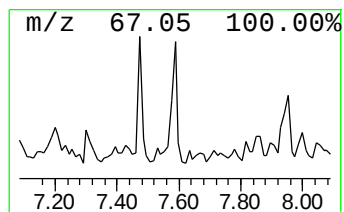
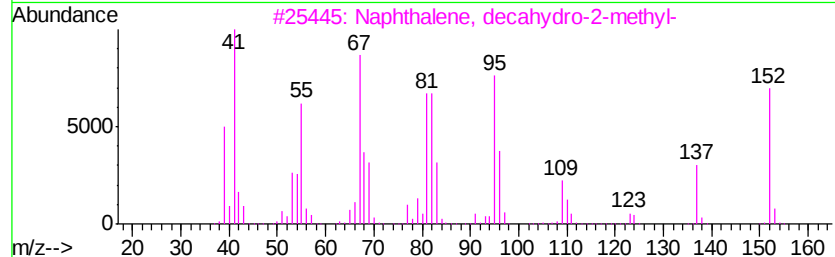
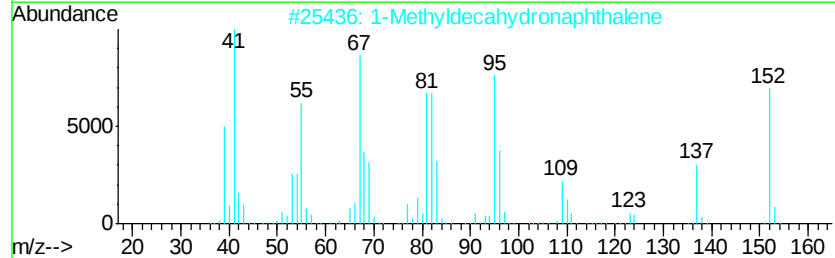
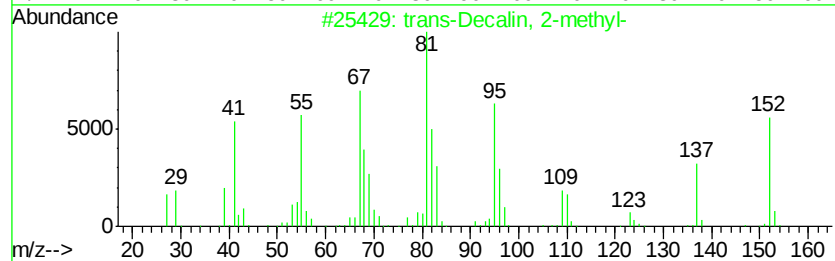
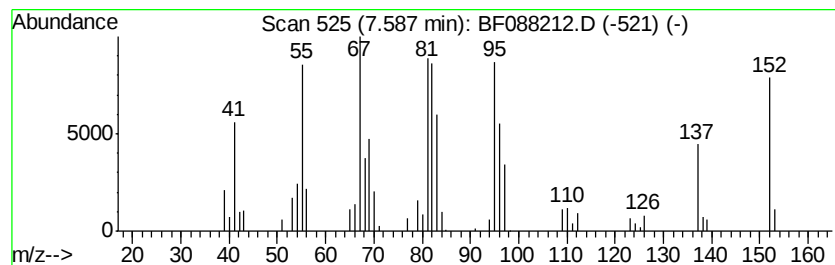
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 trans-Decalin, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	4.79 ng	152921	Naphthalene-d8	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	72
2		1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	70
3		Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	70
4		trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	64
5		Bicyclo[4.3.1]decan-10-one	152	C10H16O	020440-21-5	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
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ALS Vial : 40 Sample Multiplier: 1

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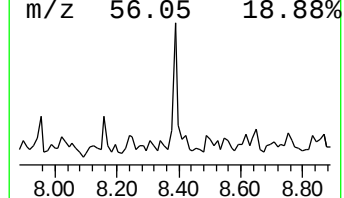
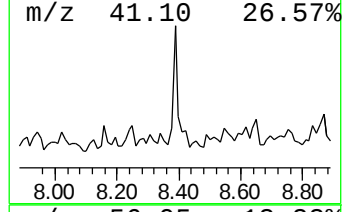
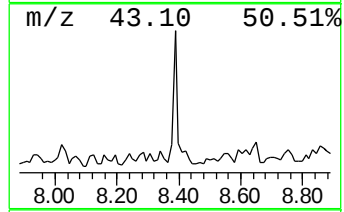
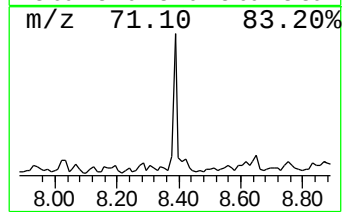
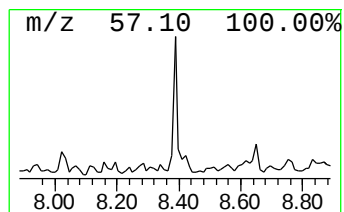
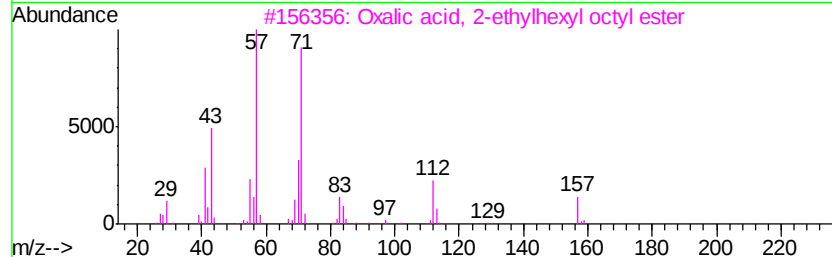
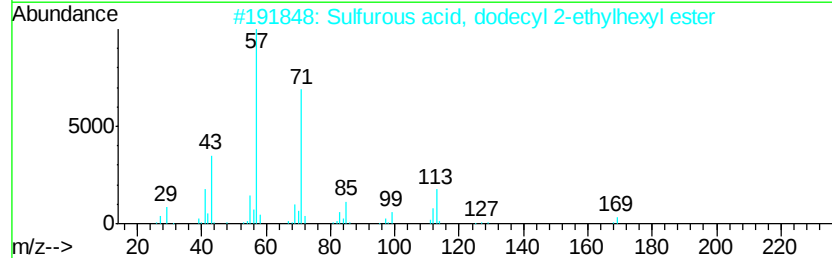
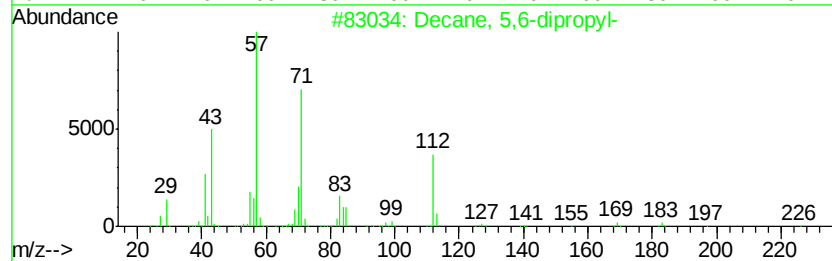
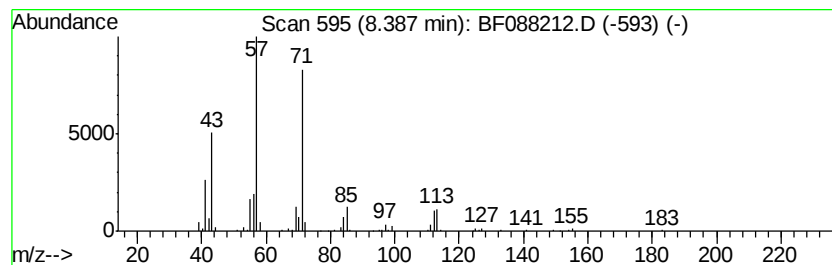
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Decane, 5,6-dipropyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.39	9.92 ng	316793	Napthalene-d8	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	78
2		Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	78
3		Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	78
4		Heptadecane, 7-methyl-	254	C18H38	020959-33-5	78
5		Decane, 4-methyl-	156	C11H24	002847-72-5	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
Sample : H3580-21
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
08-B6(8-16)C

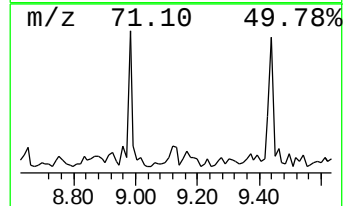
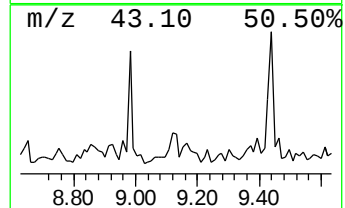
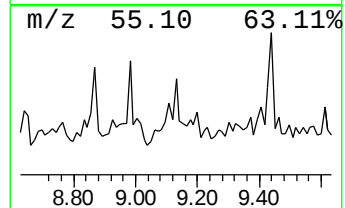
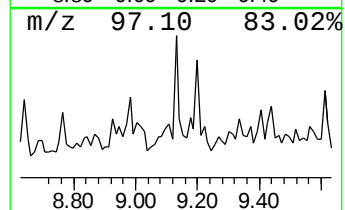
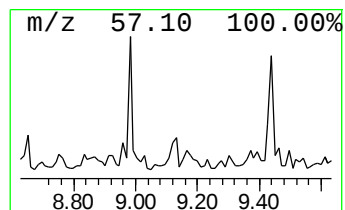
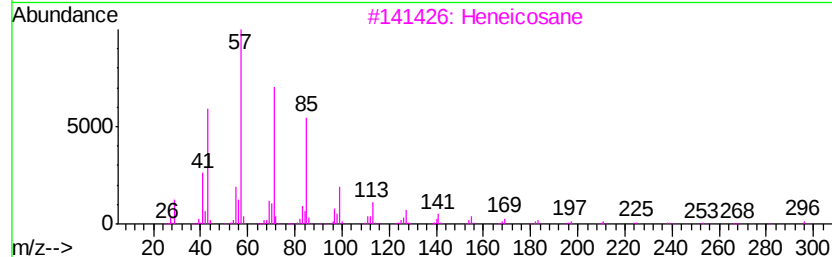
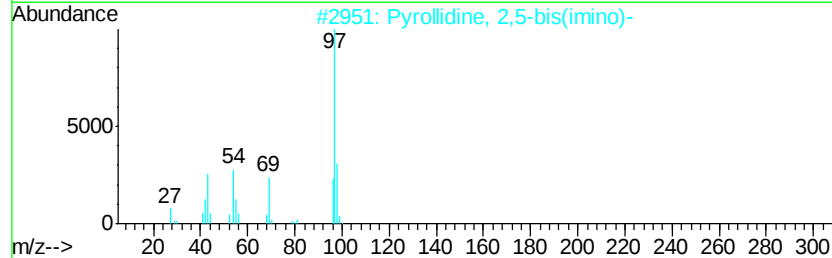
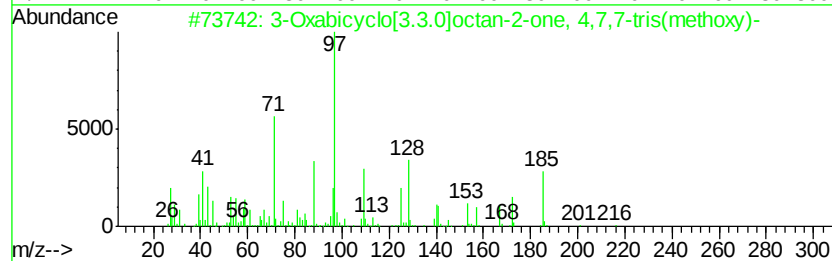
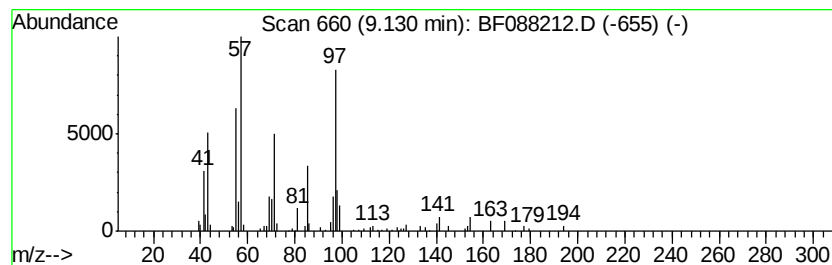
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown9.13 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.13	5.42 ng	187084	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Oxabicyclo[3.3.0]octan-2-one, ...	216	C10H16O5	1000155-61-8	47
2			Pyrollidine, 2,5-bis(imino)-	97	C4H7N3	000503-84-4	43
3			Heneicosane	296	C21H44	000629-94-7	35
4			3,5-Dimethyldodecane	198	C14H30	107770-99-0	35
5			Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
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Sample : H3580-21
Misc :
ALS Vial : 40 Sample Multiplier: 1

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ClientSampleId :
08-B6(8-16)C

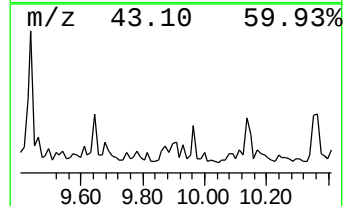
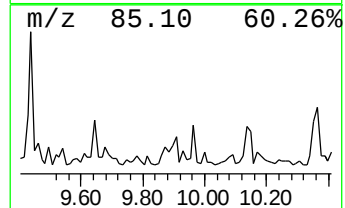
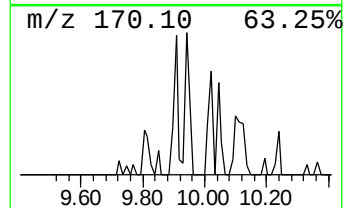
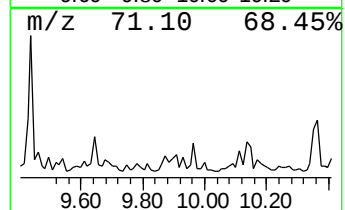
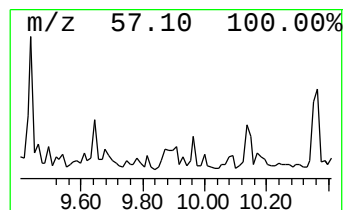
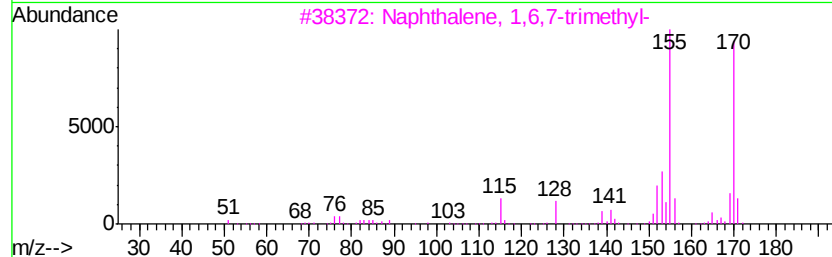
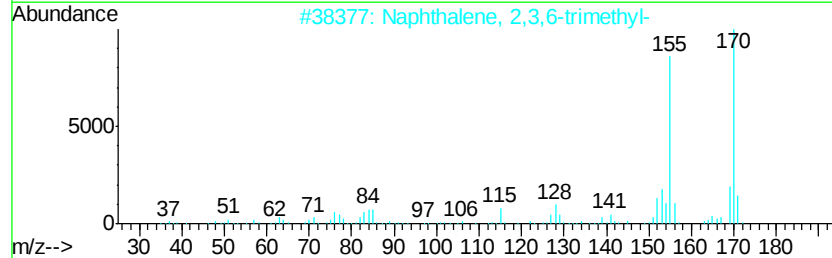
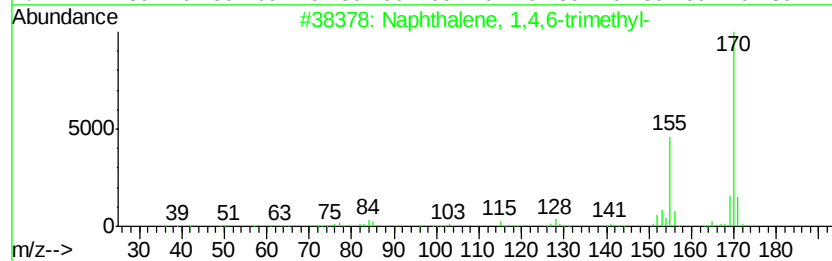
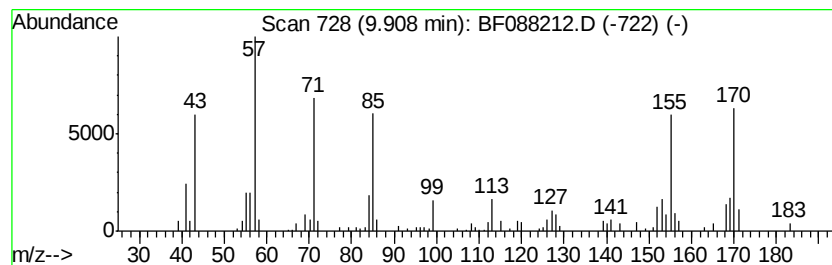
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 1,4,6-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.91	4.30 ng	148299	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	90
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	78
4			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	60
5			3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
Sample : H3580-21
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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ClientSampleId :
08-B6(8-16)C

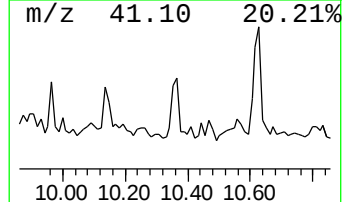
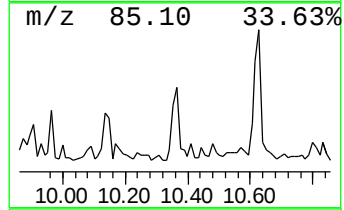
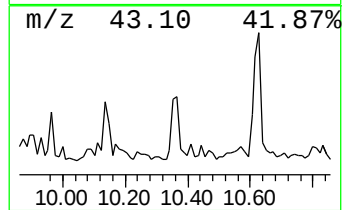
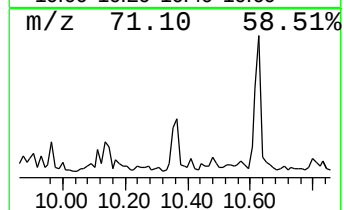
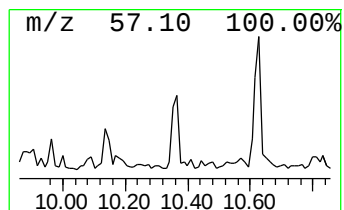
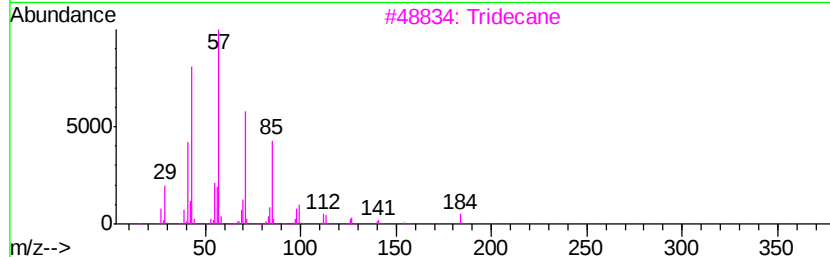
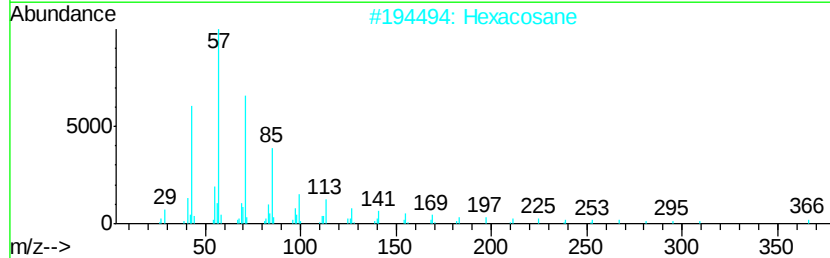
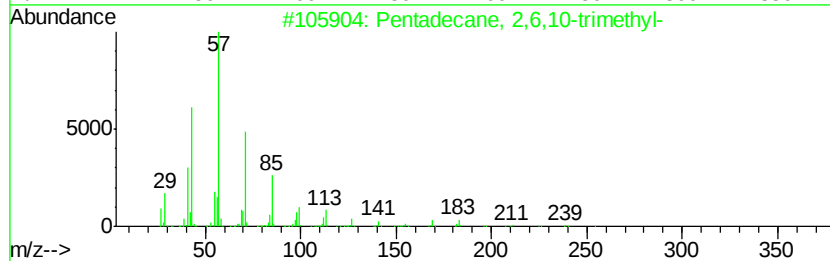
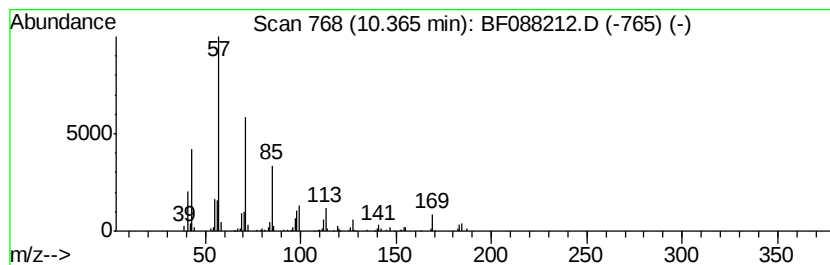
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pentadecane, 2,6,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.36	4.68 ng	161608	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	87
2			Hexacosane	366	C26H54	000630-01-3	86
3			Tridecane	184	C13H28	000629-50-5	83
4			Octadecane	254	C18H38	000593-45-3	80
5			2-Bromo dodecane	248	C12H25Br	013187-99-0	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
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Operator : UM/SJ
Sample : H3580-21
Misc :
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Instrument :
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ClientSampleId :
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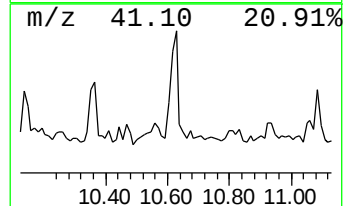
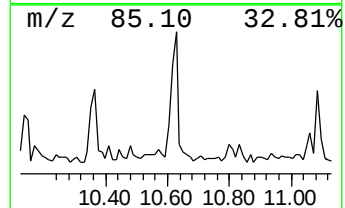
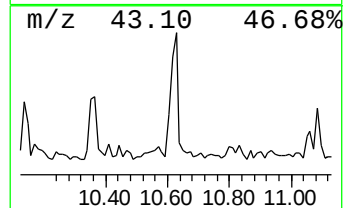
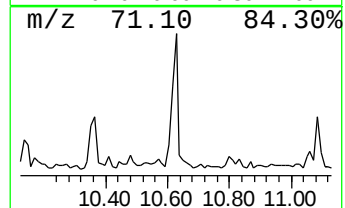
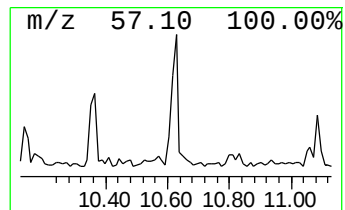
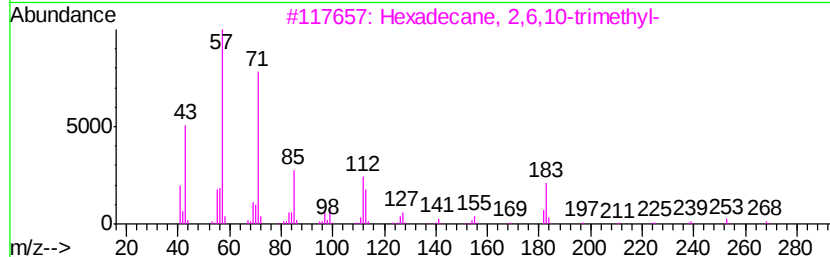
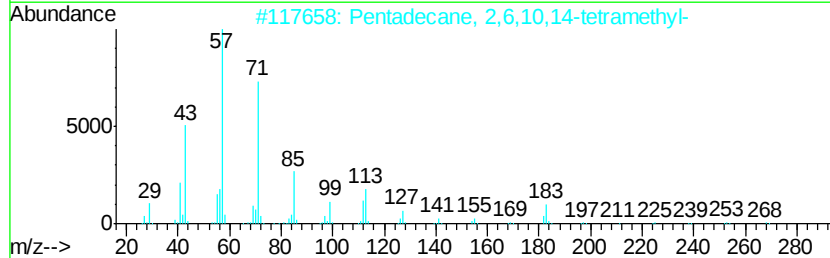
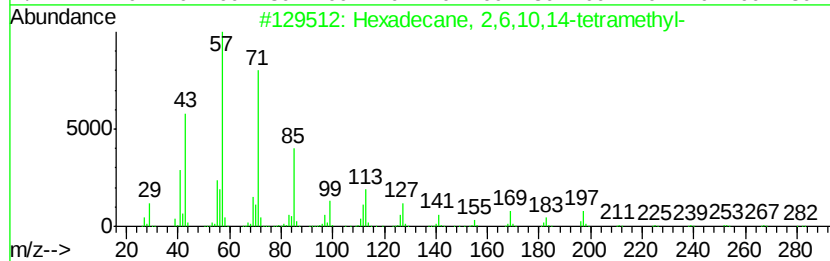
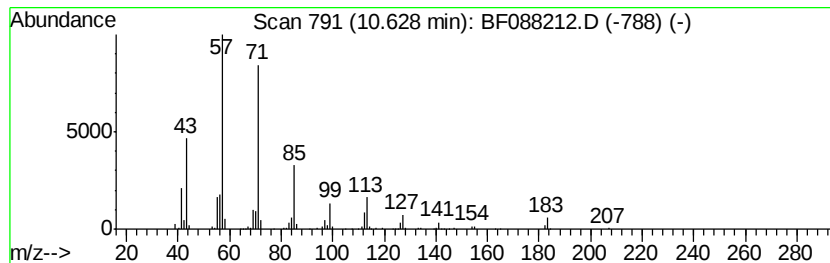
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Hexadecane, 2,6,10,14-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.63	12.26 ng	318074	Phenanthrene-d10	11.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	86
3		Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	86
4		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
5		Heptacosane	380	C27H56	000593-49-7	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
Acq On : 21 Jun 2016 11:23
Operator : UM/SJ
Sample : H3580-21
Misc :
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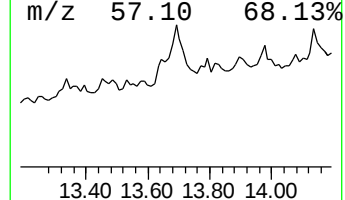
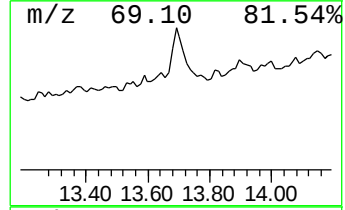
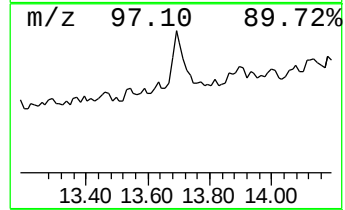
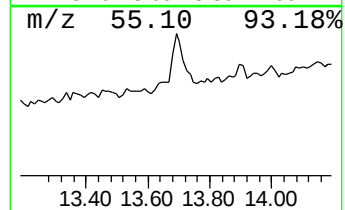
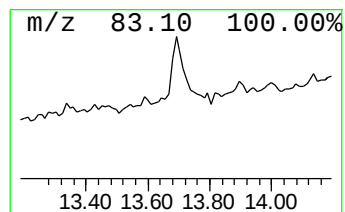
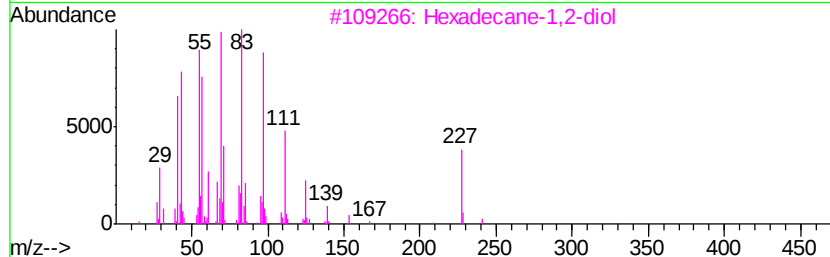
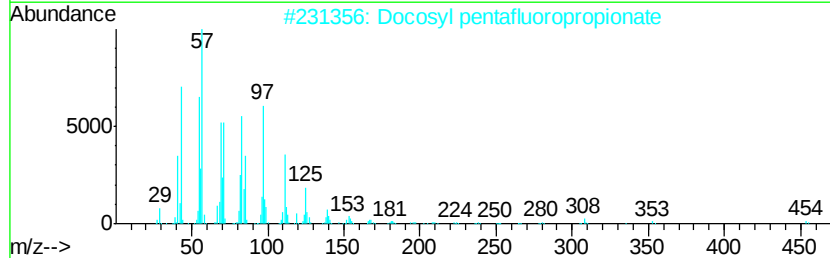
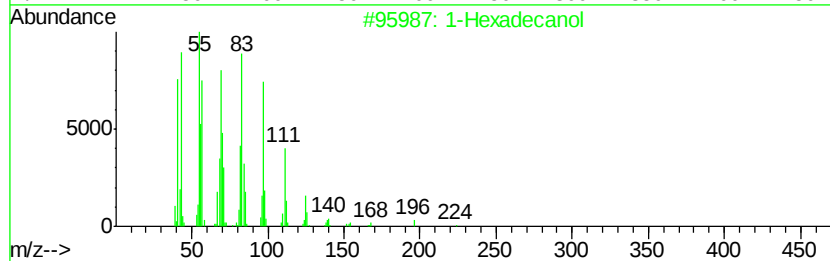
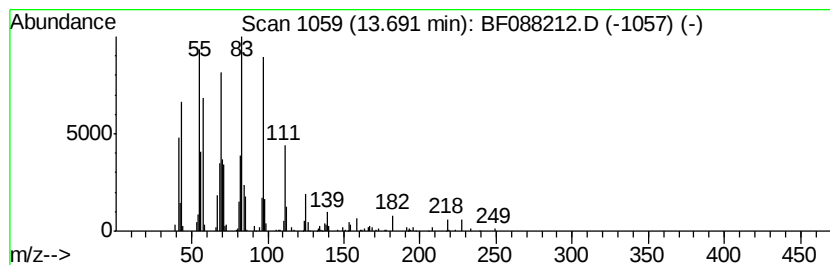
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 1-Hexadecanol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	4.08 ng	108268	Chrysene-d12	13.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Hexadecanol	242 C16H34O	036653-82-4 76
2		Docosyl pentafluoropropionate	472 C25H45F5O2	1000351-80-9 68
3		Hexadecane-1,2-diol	258 C16H34O2	006920-24-7 68
4		Trifluoroacetic acid,n-tridecyl ...	296 C15H27F3O2	053800-02-5 62
5		Cyclopentane, 1-pentyl-2-propyl-	182 C13H26	062199-51-3 60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
Data File : BF088212.D
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Misc :
ALS Vial : 40 Sample Multiplier: 1

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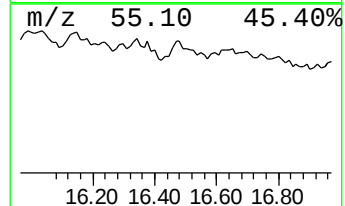
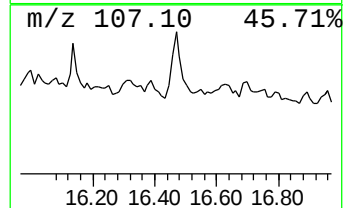
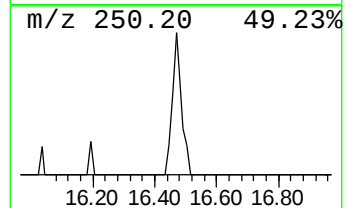
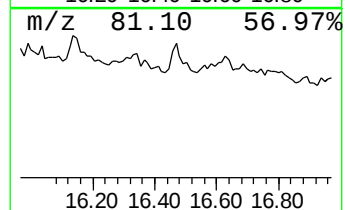
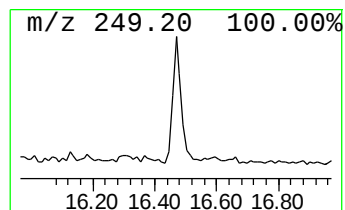
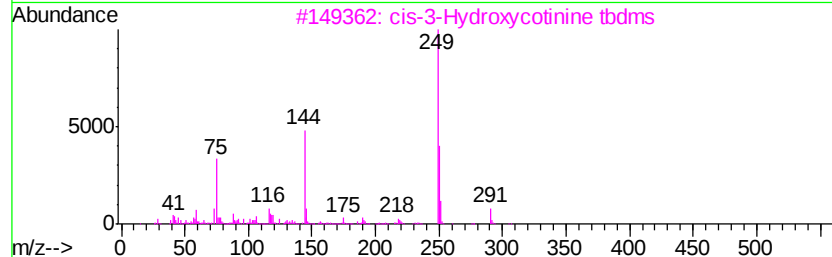
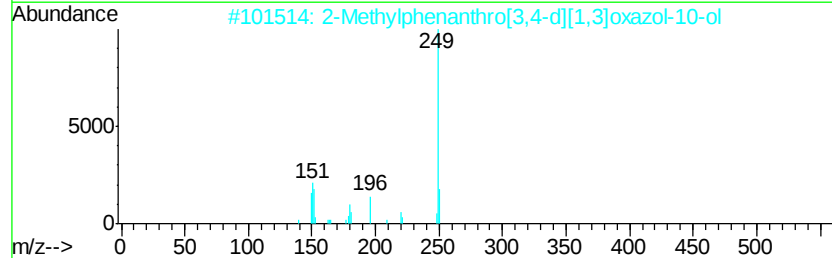
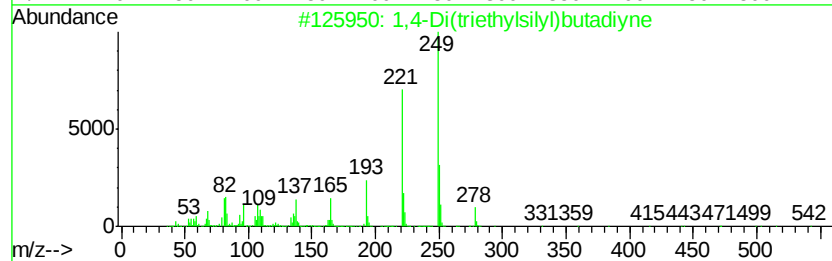
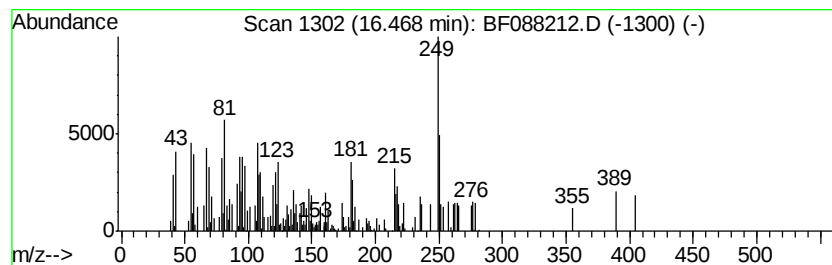
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown16.47 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	4.75 ng	129856	Perylene-d12	15.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Di(triethylsilyl)butadiyne	278	C16H30Si2	1000334-26-4	38
2			2-Methylphenanthro[3,4-d][1,3]ox...	249	C16H11N02	098033-24-0	25
3			cis-3-Hydroxycotinine tbdms	306	C16H26N2O2Si	1000331-70-7	25
4			Benzene, 1-nitro-2-hydroxy-4-(p-...	249	C12H8FN04	189214-56-0	25
5			4H-1-Benzopyran-2-carboxylic aci...	249	C12H11N05	030192-11-1	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062016\
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 ALS Vial : 40 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 08-B6(8-16)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.82	4.1 ng		100994	1	6.67	498074	20.0
Octane, 2,6-dimet...	5.91	5.4 ng		134103	1	6.67	498074	20.0
Undecane, 5,6-dim...	6.10	3.5 ng		87891	1	6.67	498074	20.0
unknown6.41	6.41	4.5 ng		110861	1	6.67	498074	20.0
unknown6.44	6.44	72.3 ng		1799440	1	6.67	498074	20.0
Nonane, 2-methyl-...	6.79	3.8 ng		95253	1	6.67	498074	20.0
unknown6.90	6.90	4.8 ng		119075	1	6.67	498074	20.0
unknown7.31	7.31	4.4 ng		108616	1	6.67	498074	20.0
trans-Decalin, 2-...	7.59	4.8 ng		152921	2	7.95	638385	20.0
Decane, 5,6-dipro...	8.39	9.9 ng		316793	2	7.95	638385	20.0
unknown9.13	9.13	5.4 ng		187084	3	9.70	689993	20.0
Naphthalene, 1,4,...	9.91	4.3 ng		148299	3	9.70	689993	20.0
Pentadecane, 2,6,...	10.36	4.7 ng		161608	3	9.70	689993	20.0
Hexadecane, 2,6,1...	10.63	12.3 ng		318074	4	11.18	518814	20.0
1-Hexadecanol	13.69	4.1 ng		108268	5	13.81	530605	20.0
unknown16.47	16.47	4.8 ng		129856	6	15.18	546213	20.0