

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
 Data File : BF094272.D
 Acq On : 7 Apr 2017 14:51
 Operator : SJ/MA
 Sample : I2528-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-6

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-BF032217.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.469	95	99	105	rVB	92925	145662	3.12%	0.171%
2	2.704	131	139	150	rBV6	52751	124941	2.68%	0.146%
3	3.028	190	194	201	rVB3	248740	417542	8.95%	0.489%
4	3.128	208	211	216	rVB	285938	370317	7.94%	0.434%
5	3.240	225	230	234	rVB3	104014	140237	3.01%	0.164%
6	3.357	244	250	253	rBV	75507	121842	2.61%	0.143%
7	3.487	268	272	279	rVB2	336545	444723	9.54%	0.521%
8	3.593	283	290	293	rBV2	213615	262738	5.63%	0.308%
9	3.922	338	346	349	rBV	231752	331737	7.11%	0.389%
10	3.987	353	357	362	rBV	392086	528934	11.34%	0.620%
11	4.175	385	389	392	rBV2	138976	193353	4.15%	0.227%
12	4.210	392	395	397	rVV2	99464	152788	3.28%	0.179%
13	4.240	397	400	405	rVV2	723614	970626	20.81%	1.137%
14	4.281	405	407	410	rVV	363364	322261	6.91%	0.378%
15	4.381	414	424	427	rVB2	3284703	4663332	100.00%	5.463%
16	4.534	445	450	455	rVB3	119386	158166	3.39%	0.185%
17	4.604	459	462	465	rBV	548158	548297	11.76%	0.642%
18	4.669	469	473	480	rVB	555050	536615	11.51%	0.629%
19	4.834	497	501	504	rBV3	106257	181729	3.90%	0.213%
20	4.934	509	518	521	rBV2	321596	519366	11.14%	0.608%
21	4.975	521	525	528	rBV2	138132	141532	3.03%	0.166%
22	5.022	530	533	536	rBV2	285515	289072	6.20%	0.339%
23	5.245	563	571	574	rBV	799084	1027344	22.03%	1.203%
24	5.275	574	576	578	rVV	158572	177133	3.80%	0.208%
25	5.298	578	580	582	rVV	371218	348840	7.48%	0.409%
26	5.334	582	586	587	rVV3	402994	580213	12.44%	0.680%
27	5.351	587	589	594	rVB	744803	670243	14.37%	0.785%
28	5.404	594	598	601	rBV2	986751	1028672	22.06%	1.205%
29	5.451	601	606	610	rVV	2694833	3591772	77.02%	4.208%
30	5.493	610	613	621	rVB	700814	744837	15.97%	0.873%
31	5.593	621	630	633	rBV	3219737	3605988	77.33%	4.224%
32	5.651	637	640	643	rVV	1884227	2052189	44.01%	2.404%
33	5.675	643	644	649	rVB	632475	432552	9.28%	0.507%
34	5.781	658	662	664	rBV	176277	200680	4.30%	0.235%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
 Data File : BF094272.D
 Acq On : 7 Apr 2017 14:51
 Operator : SJ/MA
 Sample : I2528-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-6

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-BF032217.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	5.798	664	665	668 rVV	161380	126198	2.71%	0.148%
36	5.845	668	673	676 rVV	1189420	1158045	24.83%	1.357%
37	5.881	676	679	682 rVV2	243284	254263	5.45%	0.298%
38	5.916	682	685	691 rVV	993477	1039988	22.30%	1.218%
39	6.022	696	703	706 rVV	2634797	2991553	64.15%	3.504%
40	6.051	706	708	713 rVB	1440389	1210472	25.96%	1.418%
41	6.140	718	723	726 rVV	780522	781618	16.76%	0.916%
42	6.169	726	728	733 rVV2	522277	683881	14.67%	0.801%
43	6.222	733	737	740 rVV	1819823	1870723	40.12%	2.191%
44	6.257	740	743	748 rVB2	339565	425564	9.13%	0.499%
45	6.310	748	752	756 rBV2	624693	754500	16.18%	0.884%
46	6.393	762	766	768 rBV	602315	560147	12.01%	0.656%
47	6.422	768	771	774 rVB	489273	489226	10.49%	0.573%
48	6.481	774	781	782 rBV2	2327577	2918733	62.59%	3.419%
49	6.498	782	784	792 rVB2	2013160	2304174	49.41%	2.699%
50	6.563	792	795	798 rVB	682938	592912	12.71%	0.695%
51	6.610	798	803	806 rBV3	281306	331060	7.10%	0.388%
52	6.645	806	809	812 rBV2	331688	341708	7.33%	0.400%
53	6.692	812	817	821 rVB3	168307	271449	5.82%	0.318%
54	6.751	821	827	830 rBV	1282075	1345517	28.85%	1.576%
55	6.781	830	832	837 rVB	1151959	1078109	23.12%	1.263%
56	6.845	840	843	848 rVB3	96522	118463	2.54%	0.139%
57	6.892	848	851	854 rBV	406901	482681	10.35%	0.565%
58	6.922	854	856	858 rVV	395368	382558	8.20%	0.448%
59	6.963	858	863	866 rVV2	1399693	1631480	34.99%	1.911%
60	6.992	866	868	872 rVV3	252090	406908	8.73%	0.477%
61	7.045	872	877	880 rVV3	1989584	2691844	57.72%	3.153%
62	7.092	880	885	889 rVB3	881495	1059836	22.73%	1.242%
63	7.151	889	895	900 rVB5	550742	1069420	22.93%	1.253%
64	7.210	900	905	909 rVB	521442	579434	12.43%	0.679%
65	7.316	918	923	925 rVV3	387295	591762	12.69%	0.693%
66	7.345	925	928	930 rVV	1949743	2407802	51.63%	2.821%
67	7.375	930	933	938 rVV3	2291003	3211224	68.86%	3.762%
68	7.416	938	940	945 rVV	555058	567404	12.17%	0.665%
69	7.475	945	950	953 rVV2	415330	556906	11.94%	0.652%
70	7.522	955	958	961 rVV4	113428	144000	3.09%	0.169%
71	7.587	965	969	972 rVB3	338512	386584	8.29%	0.453%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
 Data File : BF094272.D
 Acq On : 7 Apr 2017 14:51
 Operator : SJ/MA
 Sample : I2528-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-6

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-BF032217.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	7.692	981	987	990	rBV2	412997	656049	14.07%	0.769%
73	7.728	990	993	996	rVV3	209429	252678	5.42%	0.296%
74	7.763	996	999	1003	rVB4	155776	209096	4.48%	0.245%
75	7.804	1003	1006	1007	rBV3	193334	201898	4.33%	0.237%
76	7.828	1007	1010	1012	rVV	592896	514953	11.04%	0.603%
77	7.851	1012	1014	1019	rVB4	205867	245106	5.26%	0.287%
78	7.904	1019	1023	1025	rBV2	337398	433975	9.31%	0.508%
79	7.922	1025	1026	1029	rVB	298553	236395	5.07%	0.277%
80	7.963	1029	1033	1036	rBV	138403	164069	3.52%	0.192%
81	8.069	1049	1051	1053	rVB	179377	127260	2.73%	0.149%
82	8.116	1053	1059	1062	rBV	549365	640425	13.73%	0.750%
83	8.216	1069	1076	1083	rVB	1467725	1604971	34.42%	1.880%
84	8.334	1089	1096	1099	rBV	813542	874584	18.75%	1.025%
85	8.669	1145	1153	1157	rVB	3429721	3665709	78.61%	4.294%
86	8.810	1174	1177	1181	rVB	222751	223767	4.80%	0.262%
87	8.981	1203	1206	1212	rVB2	140492	214650	4.60%	0.251%
88	9.075	1215	1222	1225	rBV	159874	206890	4.44%	0.242%
89	9.163	1233	1237	1240	rVV	344137	349073	7.49%	0.409%
90	9.210	1240	1245	1252	rVB3	93084	203662	4.37%	0.239%
91	9.486	1284	1292	1295	rBV2	1397720	1653825	35.46%	1.937%
92	10.080	1389	1393	1398	rVB	120964	116713	2.50%	0.137%
93	10.457	1451	1457	1463	rBV	1969278	2092169	44.86%	2.451%
94	10.663	1488	1492	1494	rBV	123824	129486	2.78%	0.152%
95	11.304	1596	1601	1604	rVV	1296383	1369642	29.37%	1.604%
96	11.333	1604	1606	1611	rVB	182782	151523	3.25%	0.178%
97	13.286	1932	1938	1942	rBV	2648856	2850757	61.13%	3.340%
98	14.451	2132	2136	2144	rBV	213817	317104	6.80%	0.371%
99	14.557	2149	2154	2158	rBV	968440	1012295	21.71%	1.186%
100	16.186	2426	2431	2438	rVB	680952	800571	17.17%	0.938%

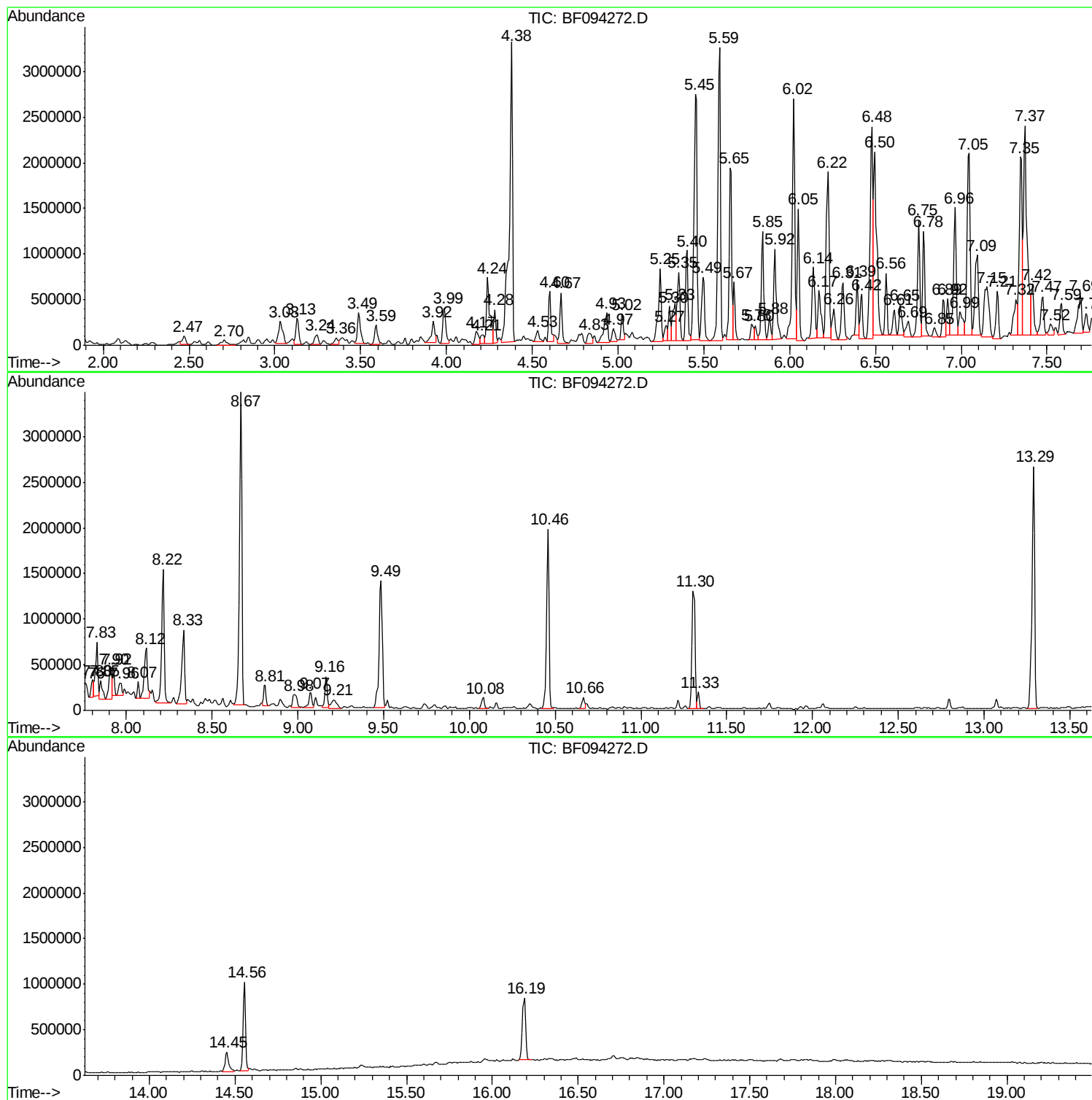
Sum of corrected areas: 85363714

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
TS-6

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
TS-6

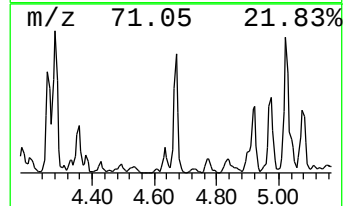
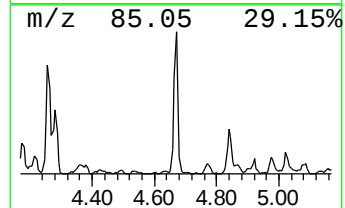
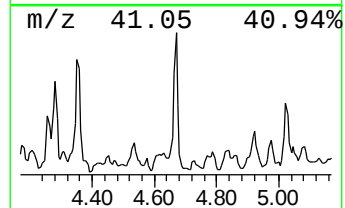
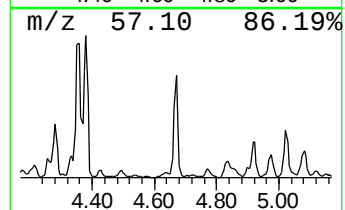
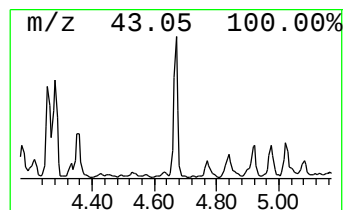
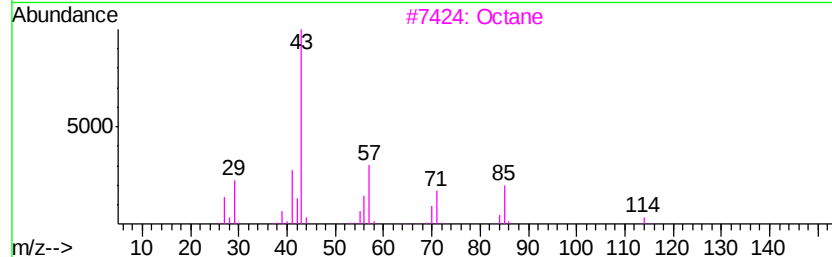
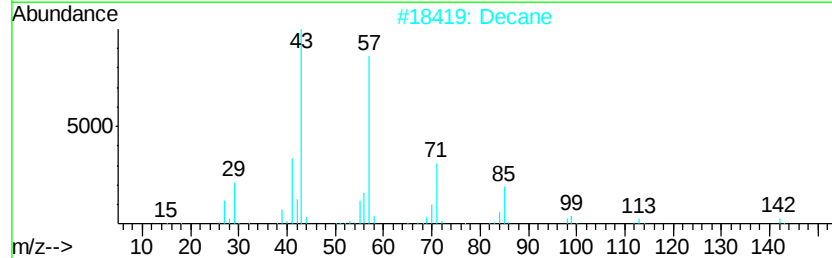
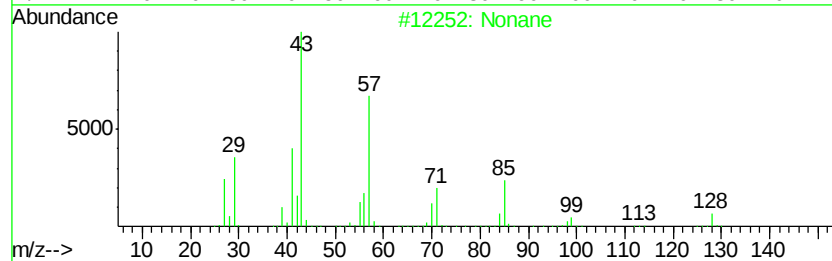
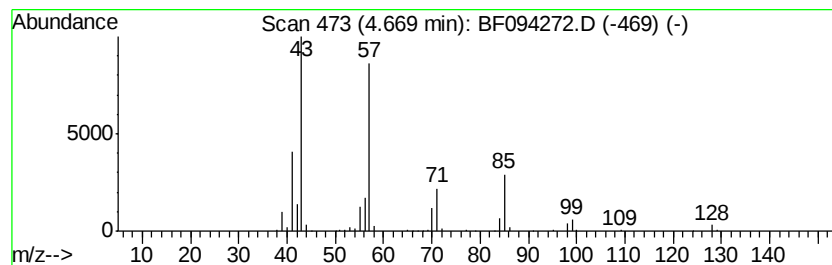
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Nonane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.67	9.27 ng	536615	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane	128	C9H20	000111-84-2	90
2		Decane	142	C10H22	000124-18-5	78
3		Octane	114	C8H18	000111-65-9	64
4		Hexadecane	226	C16H34	000544-76-3	64
5		Heptane, 4-ethyl-	128	C9H20	002216-32-2	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

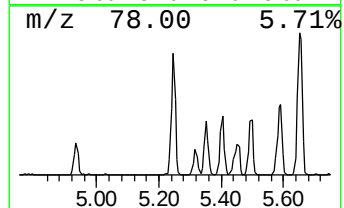
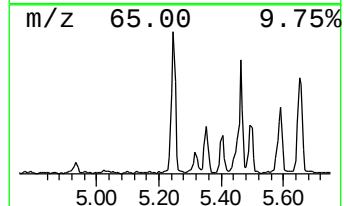
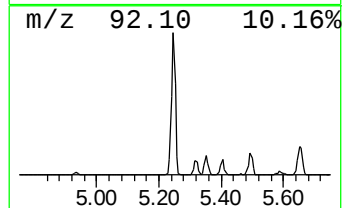
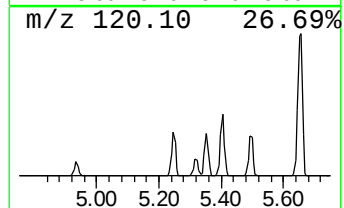
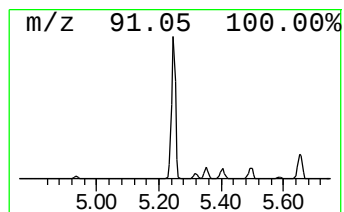
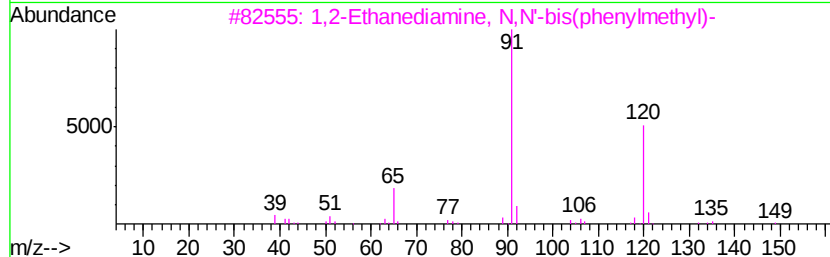
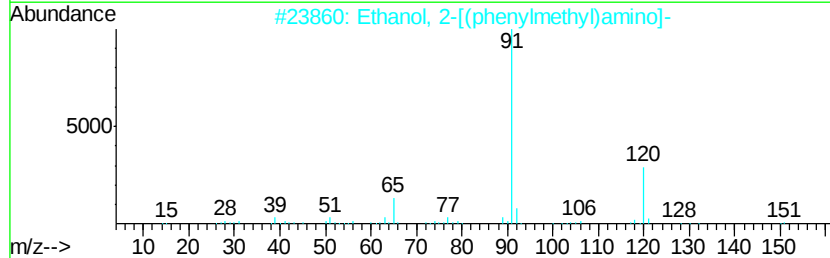
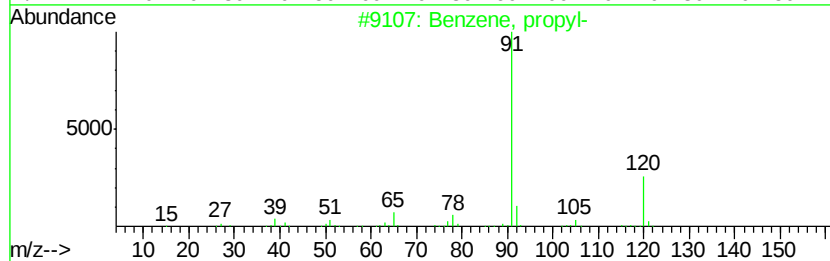
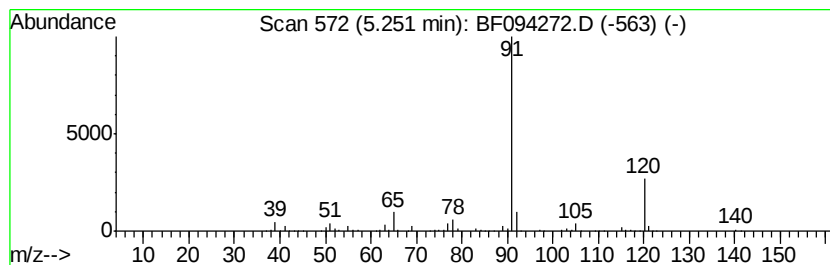
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	17.74 ng	1027340	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	90
2		Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
3		1,2-Ethanediamine, N,N'-bis(phen...)	240	C16H20N2	000140-28-3	64
4		Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	56
5		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

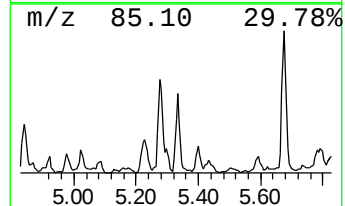
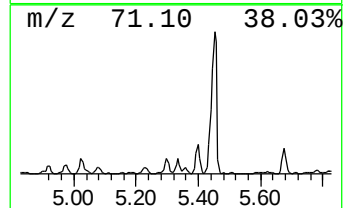
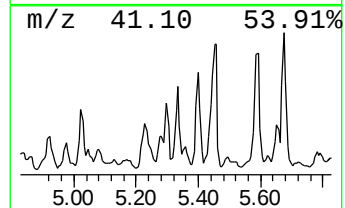
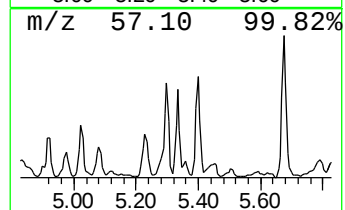
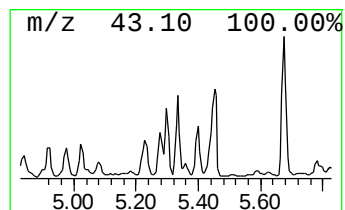
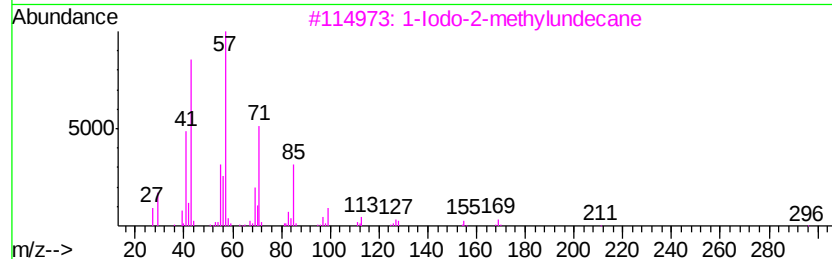
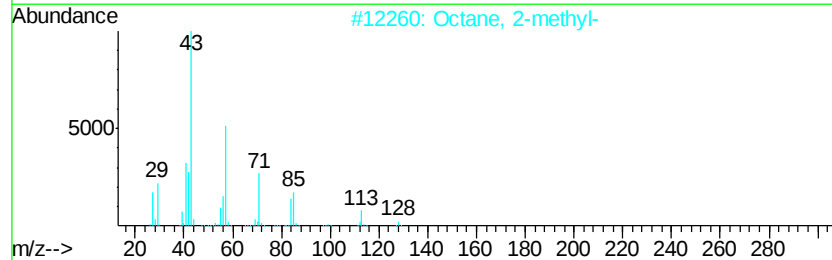
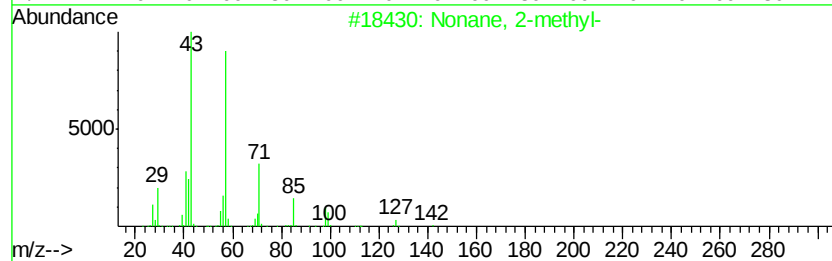
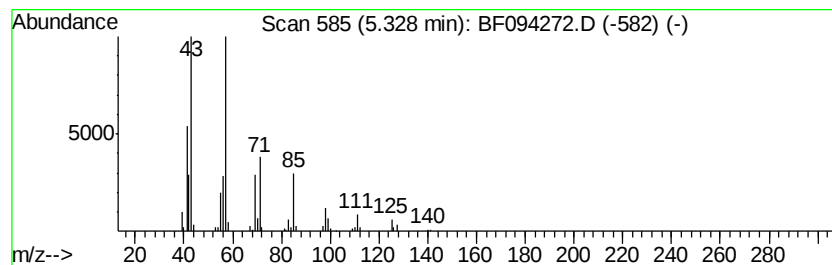
Instrument :
BNA_F
ClientSampleId :
TS-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonane, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
5.33	10.02 ng	580213	1,4-Dichlorobenzene-d4	5.85		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-	142	C10H22	000871-83-0	74	
2	Octane, 2-methyl-	128	C9H20	003221-61-2	59	
3	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	59	
4	Decane, 2-methyl-	156	C11H24	006975-98-0	53	
5	1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	53	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

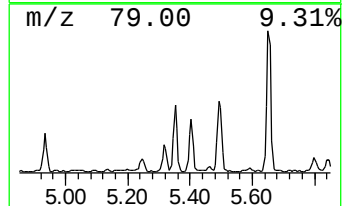
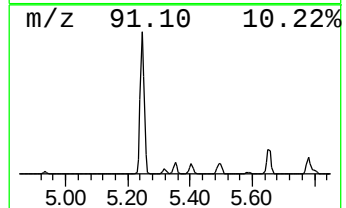
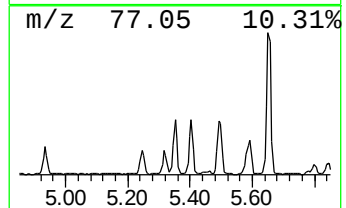
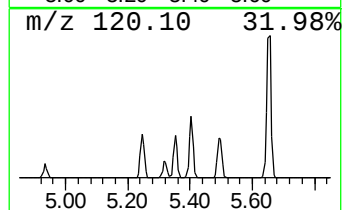
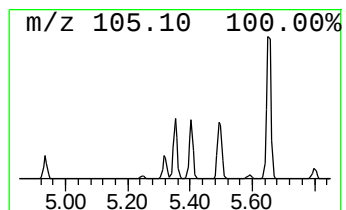
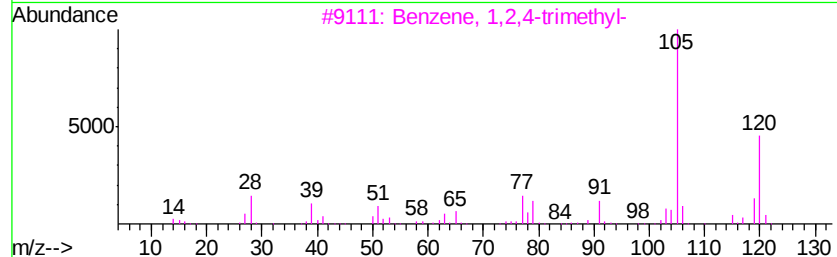
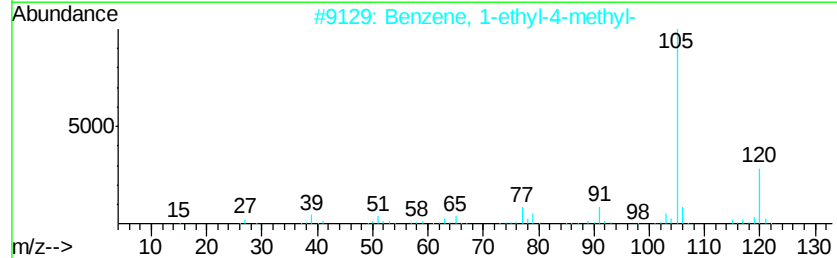
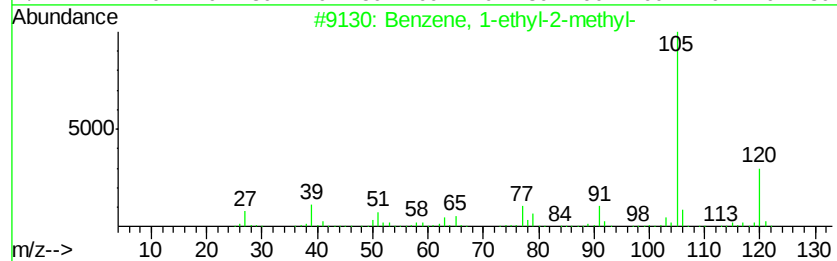
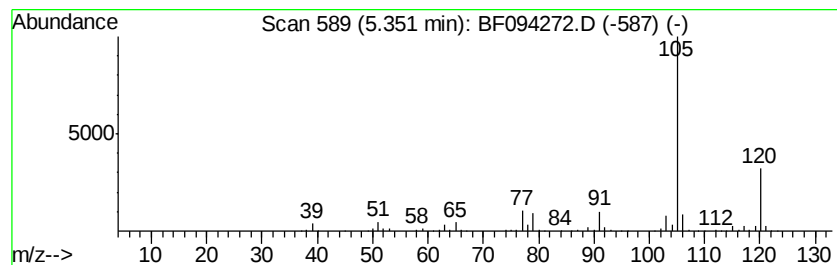
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.35	11.58 ng	670243	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

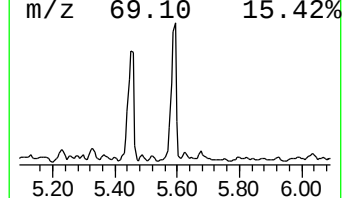
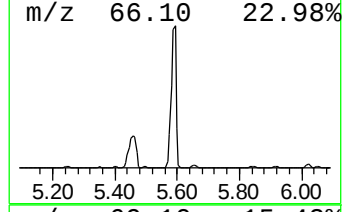
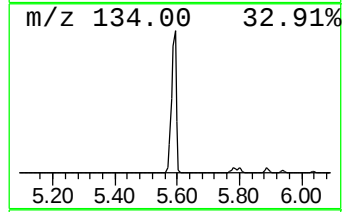
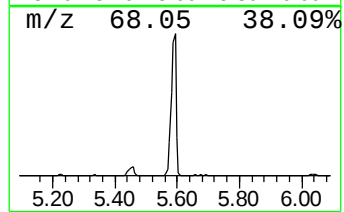
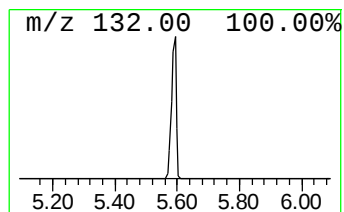
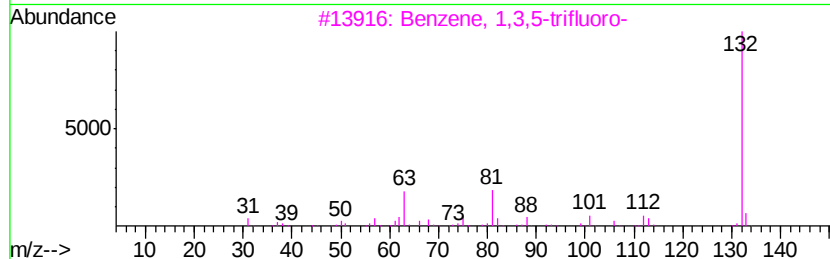
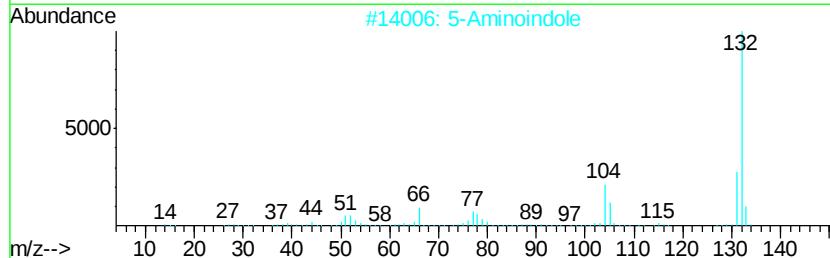
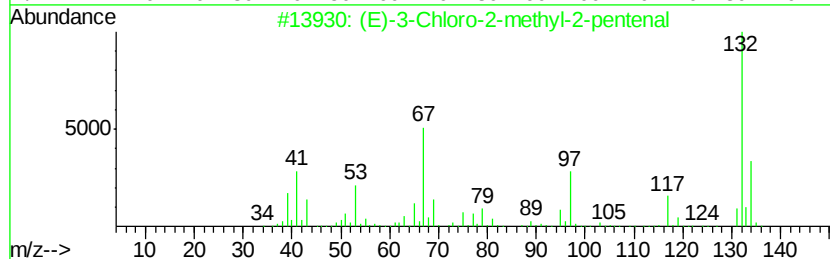
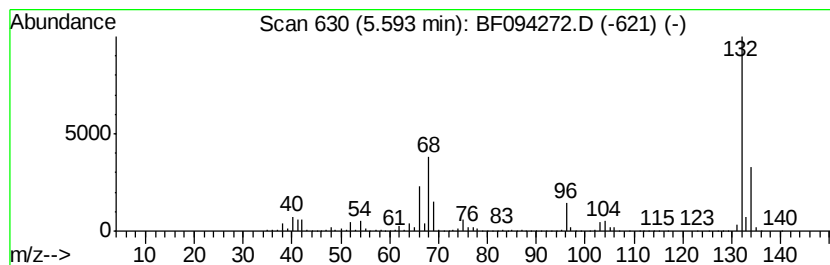
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown5.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.59	62.28 ng	3605990	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

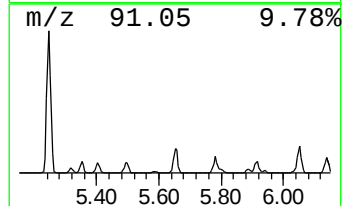
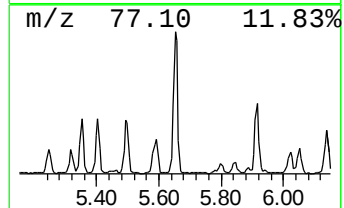
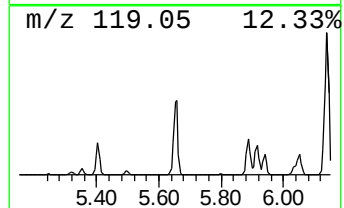
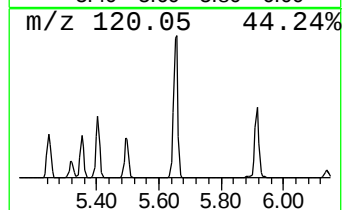
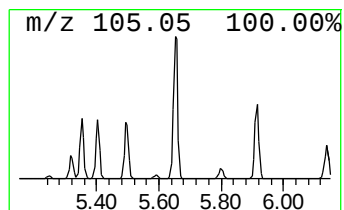
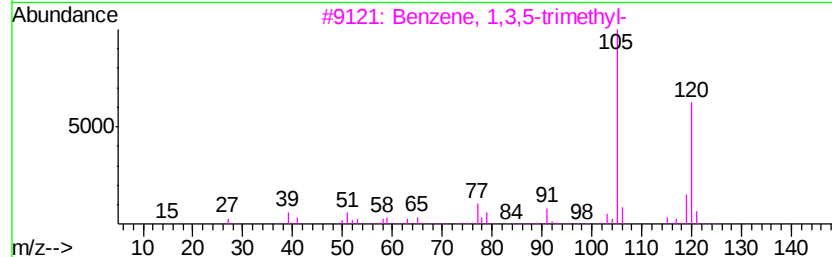
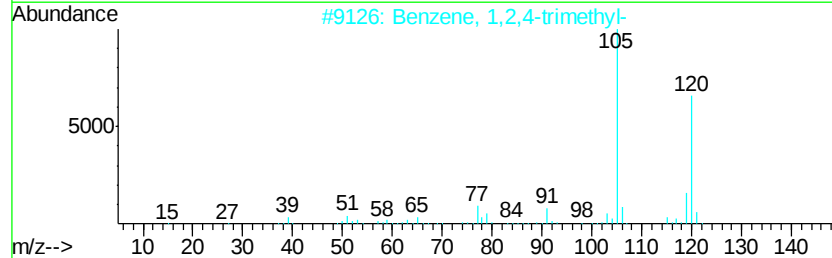
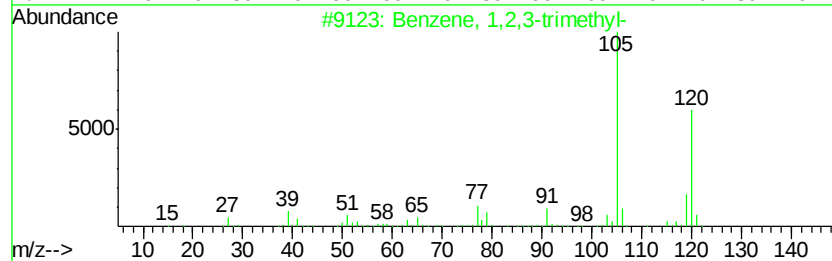
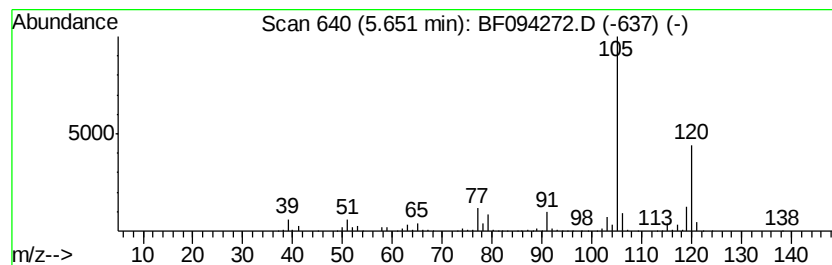
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.65	35.44 ng	2052190	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
4		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

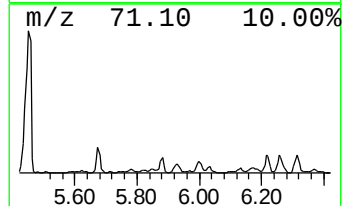
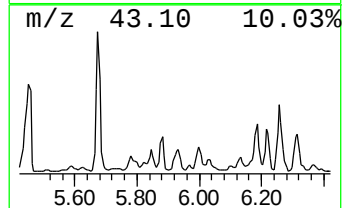
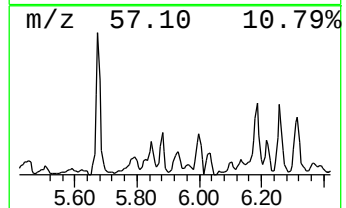
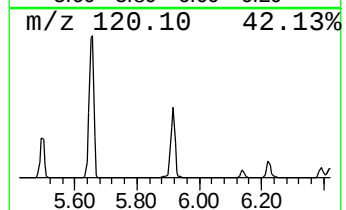
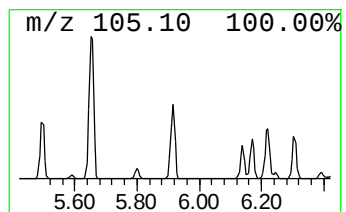
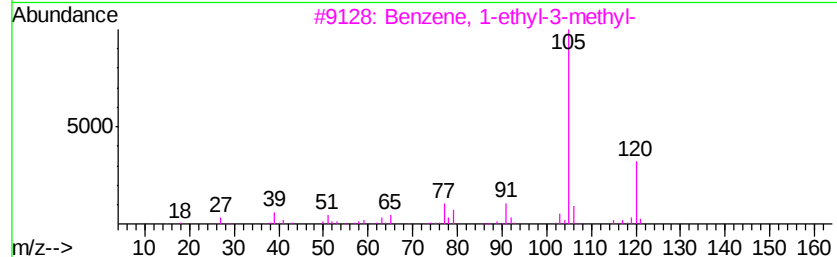
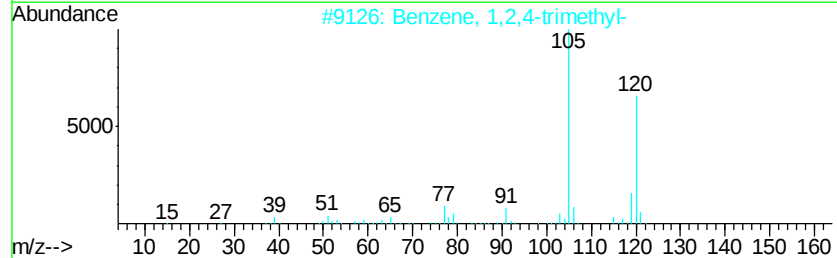
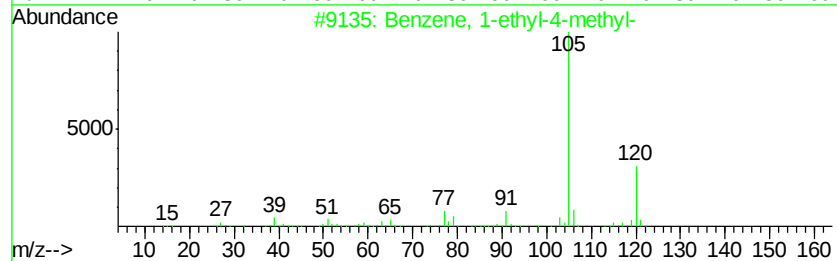
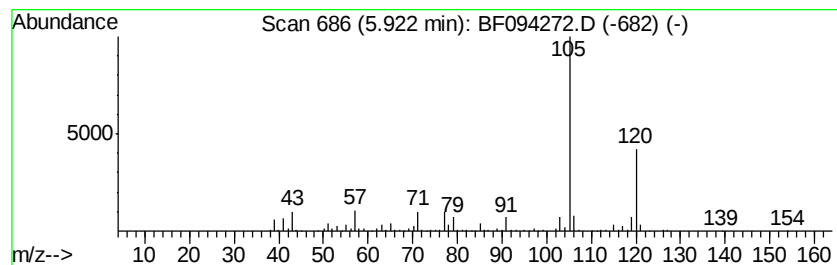
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.92	17.96 ng	1039990	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

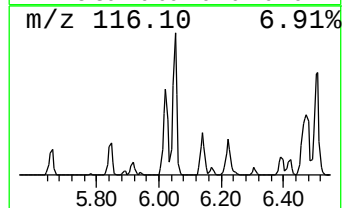
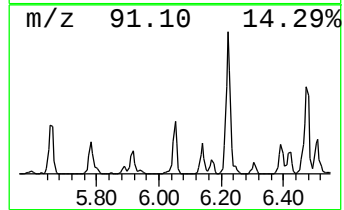
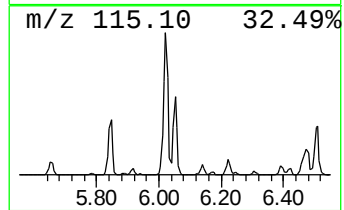
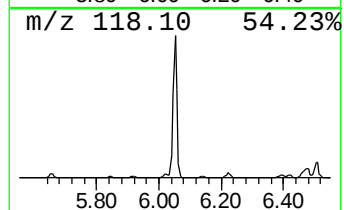
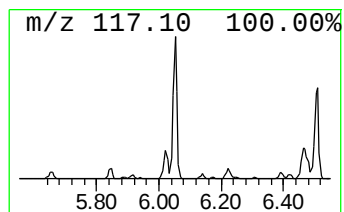
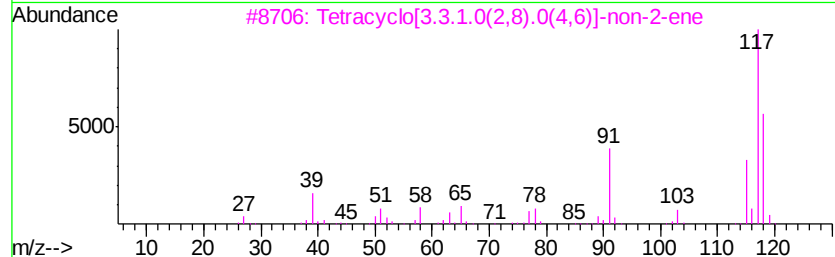
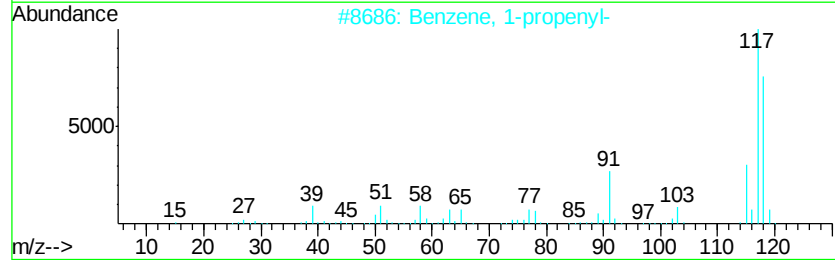
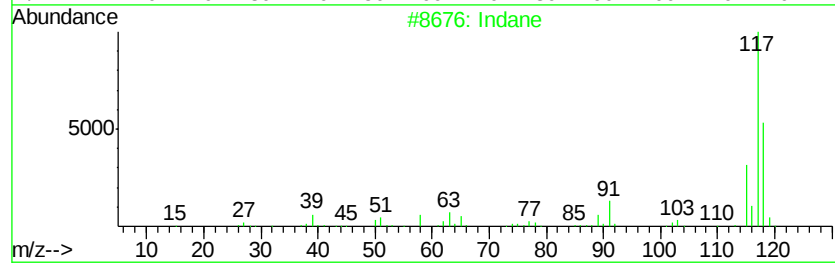
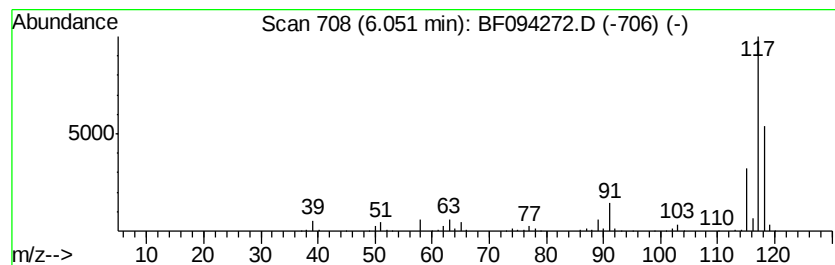
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.05	20.91 ng	1210470	1,4-Dichlorobenzene-d4	5.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	72
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	72
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	72
4			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
5			7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

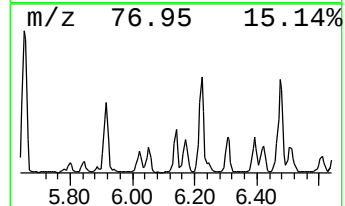
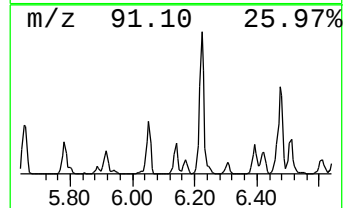
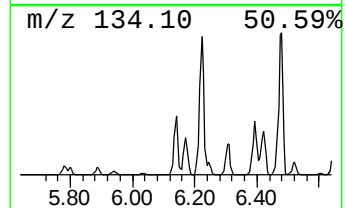
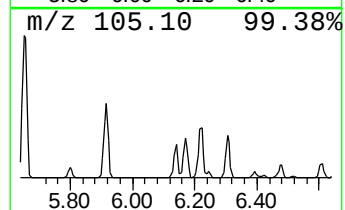
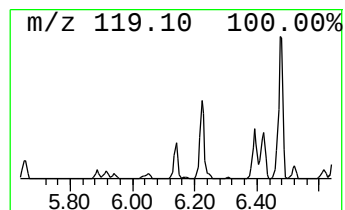
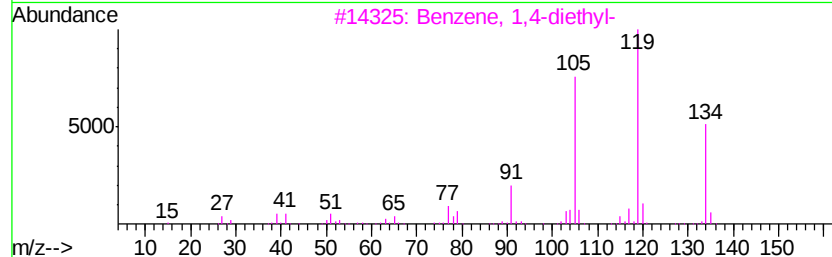
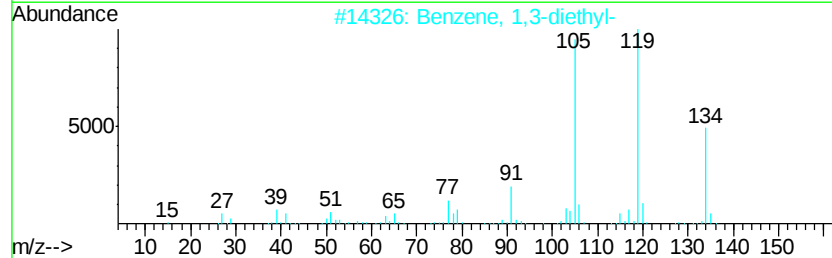
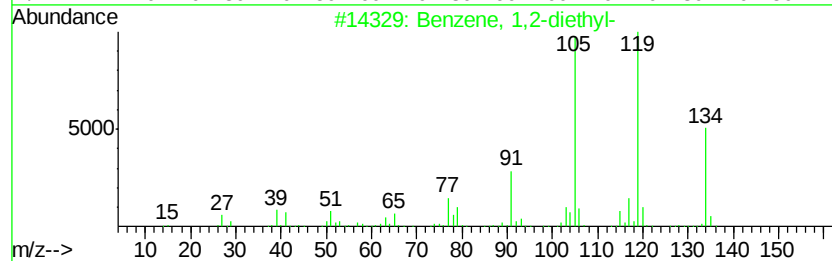
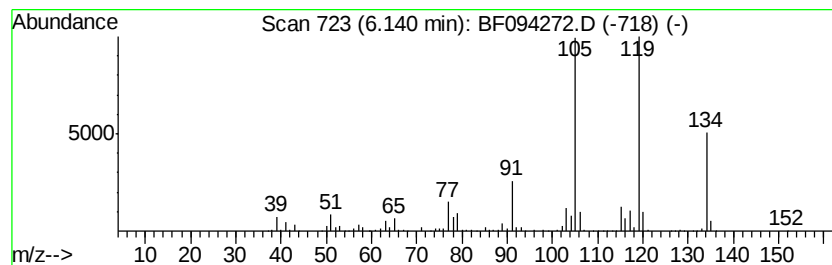
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1,2-diethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	13.50 ng	781618	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

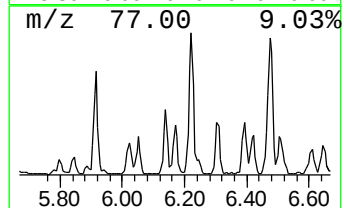
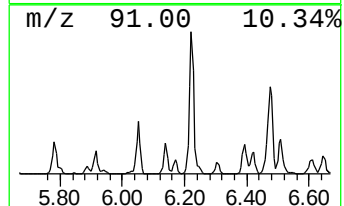
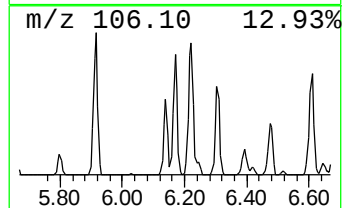
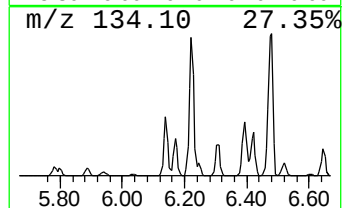
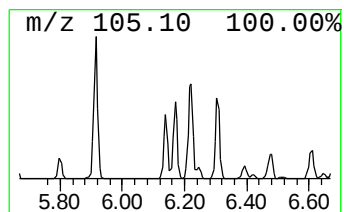
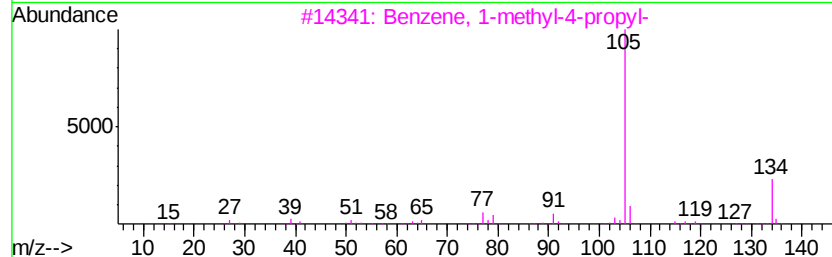
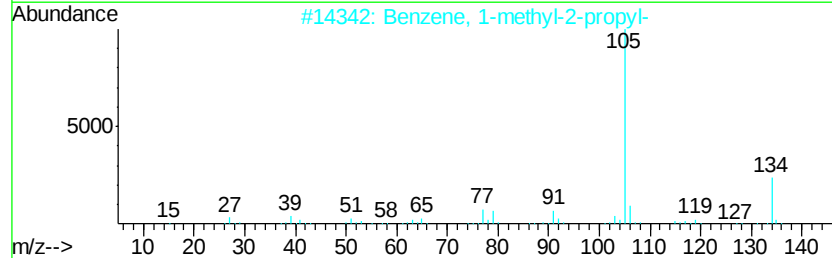
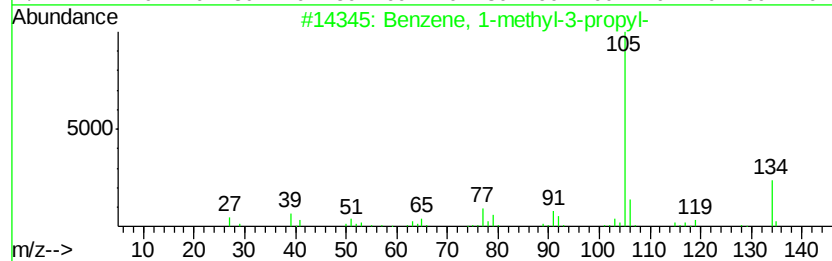
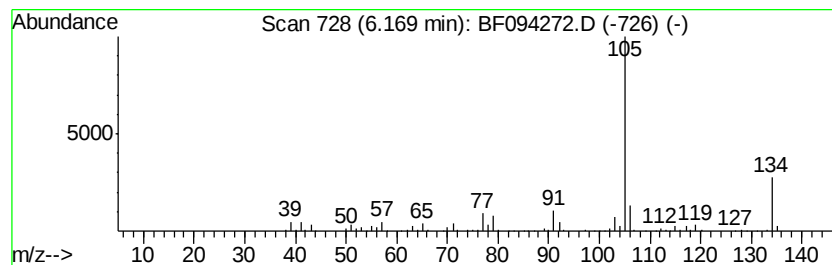
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 1-methyl-3-propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.17	11.81 ng	683881	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	94
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
TS-6

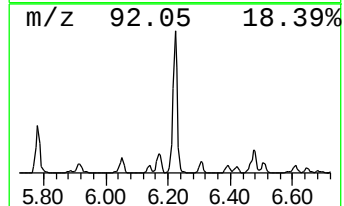
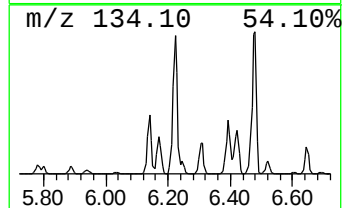
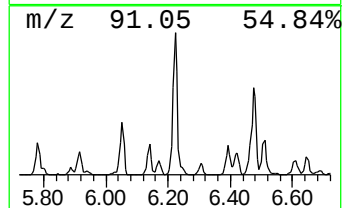
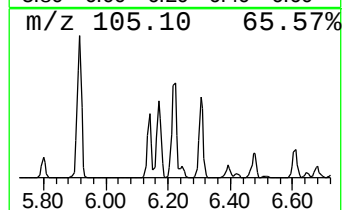
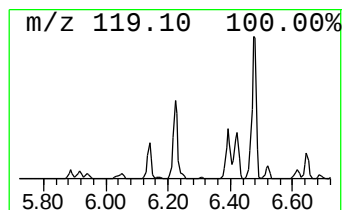
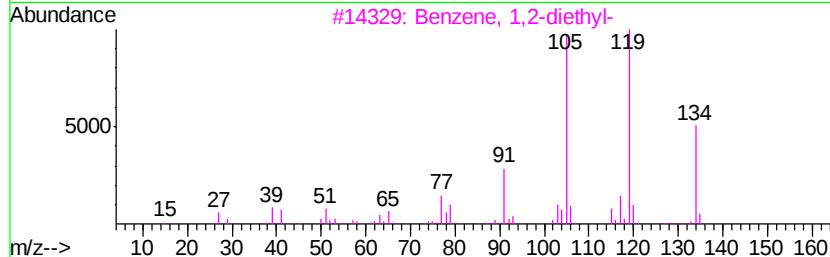
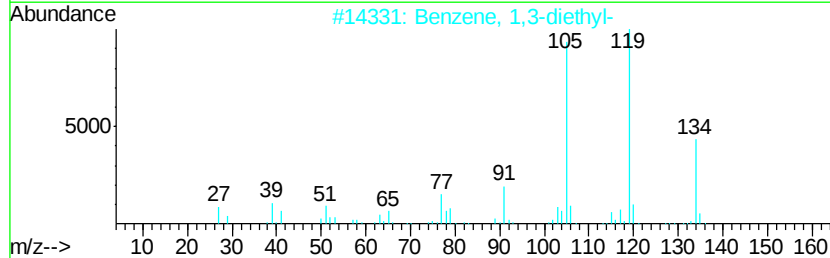
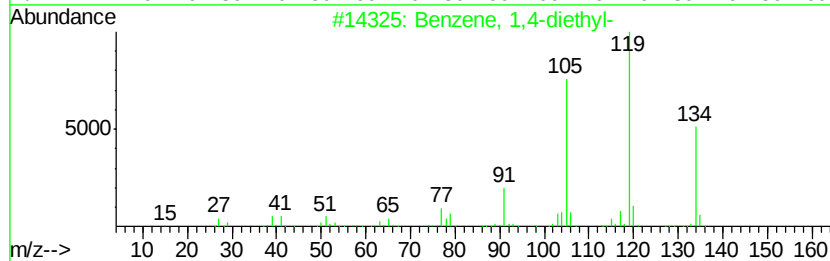
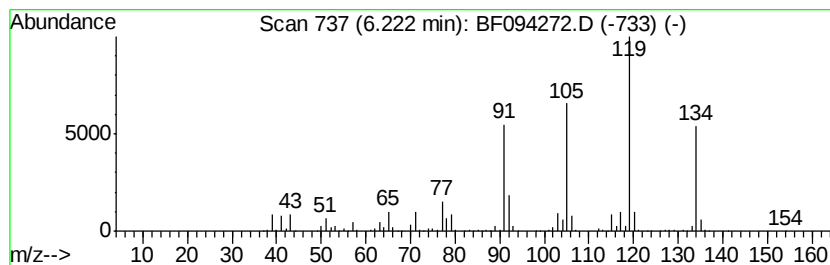
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzene, 1,4-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.22	32.31 ng	1870720	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

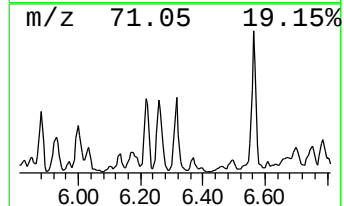
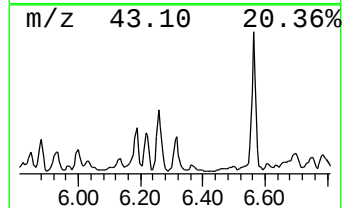
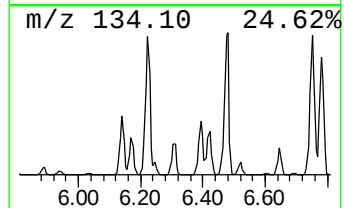
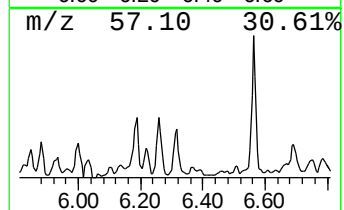
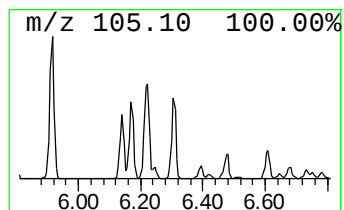
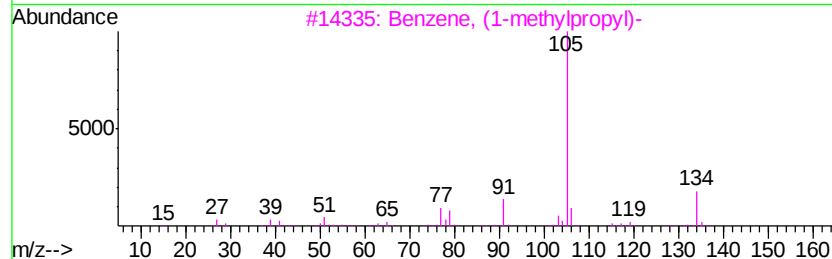
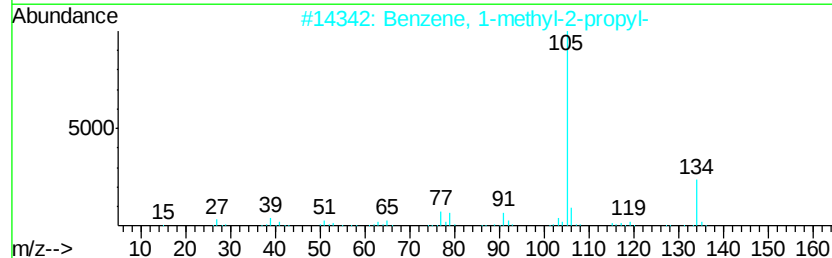
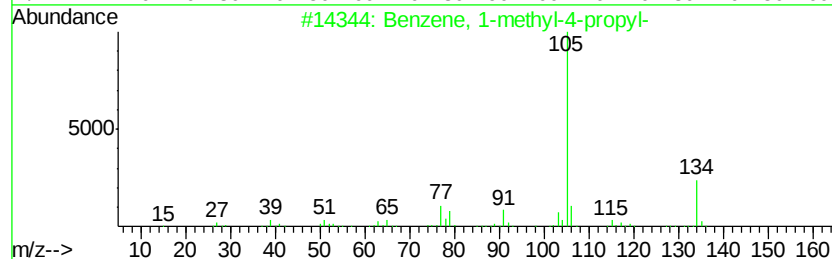
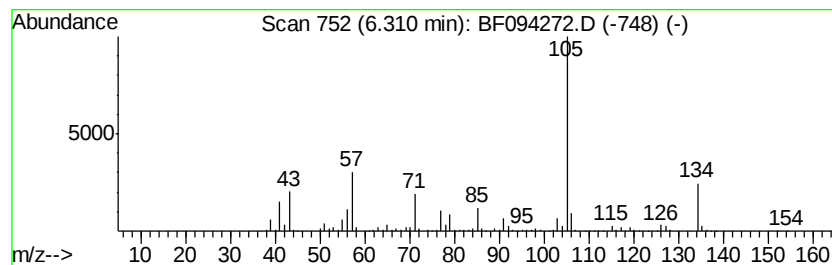
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzene, 1-methyl-4-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	13.03 ng	754500	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
2		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
3		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	60
4		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	60
5		2-Tolylloxirane	134	C9H10O	002783-26-8	45



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

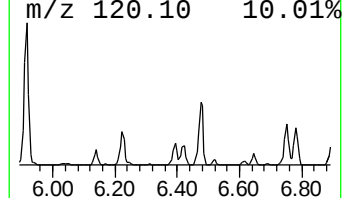
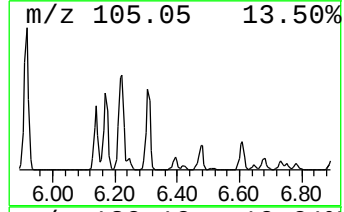
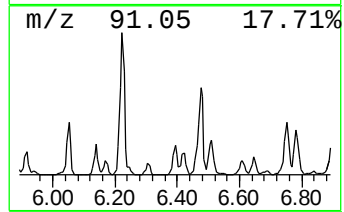
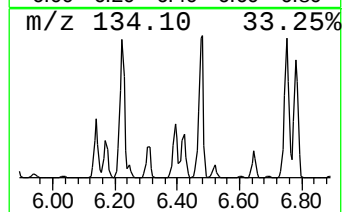
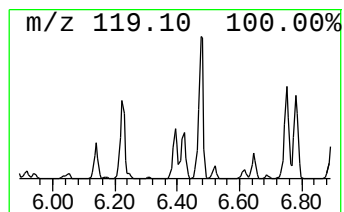
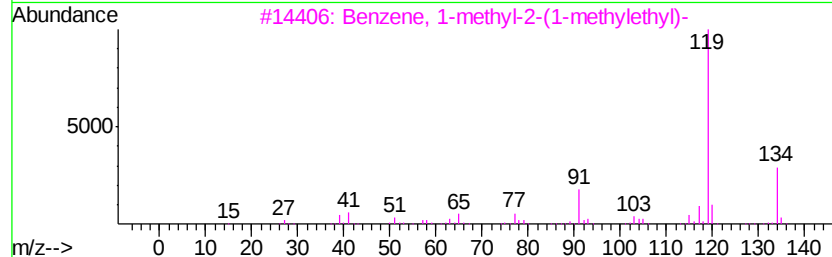
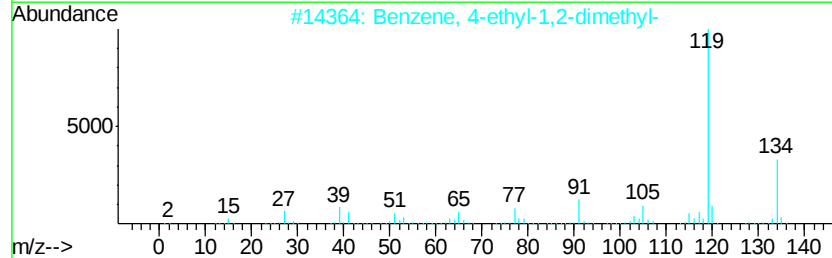
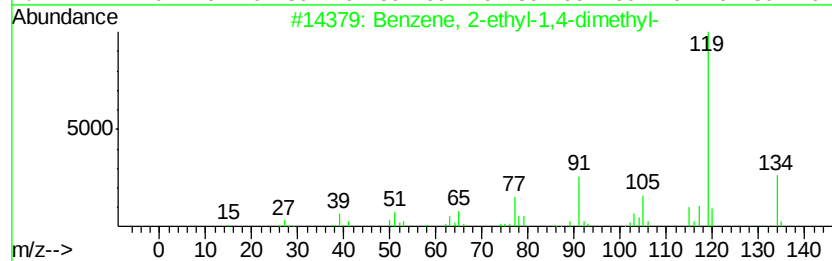
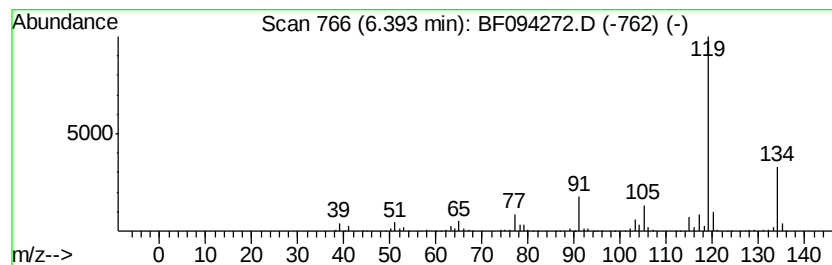
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	9.67 ng	560147	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-methyl-2-(1-methylethyl)-	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

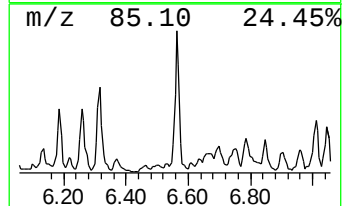
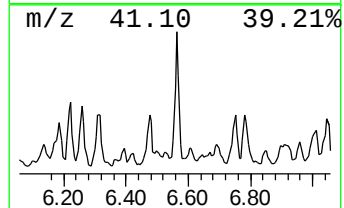
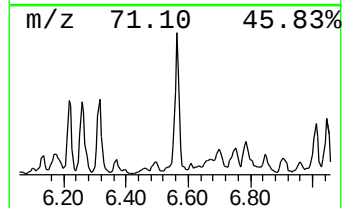
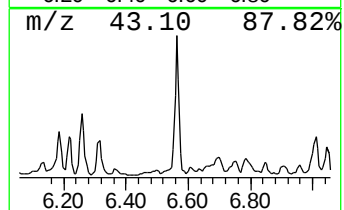
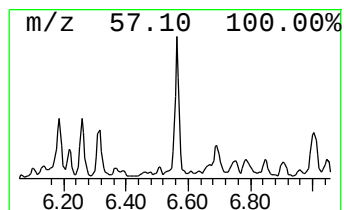
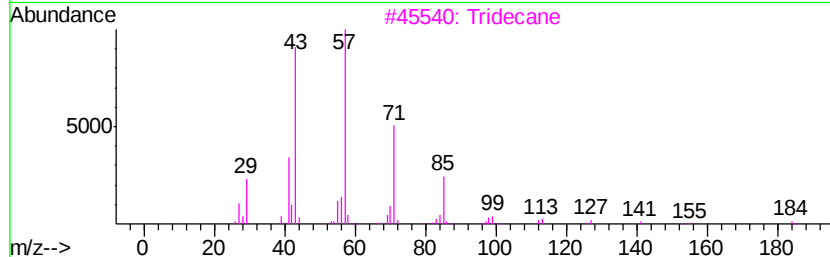
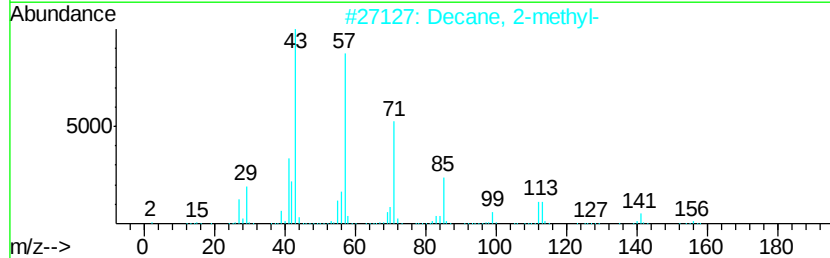
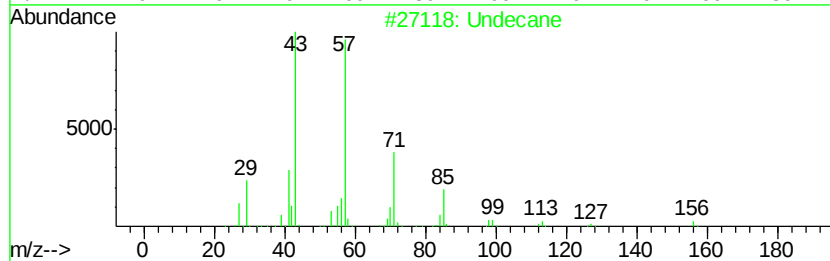
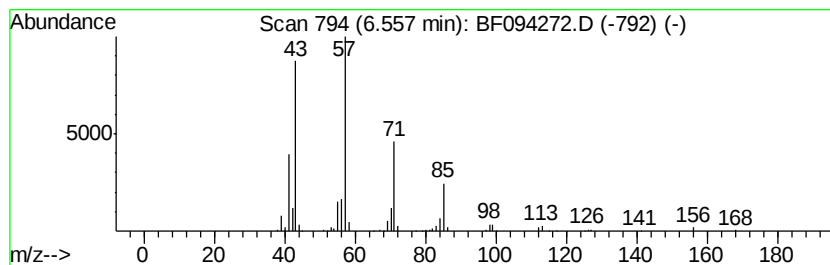
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Undecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	10.24 ng	592912	1,4-Dichlorobenzene-d4	5.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	94
2	Decane, 2-methyl-			156	C11H24	006975-98-0	80
3	Tridecane			184	C13H28	000629-50-5	78
4	Tetradecane			198	C14H30	000629-59-4	64
5	Octane, 3-ethyl-2,7-dimethyl-			170	C12H26	062183-55-5	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

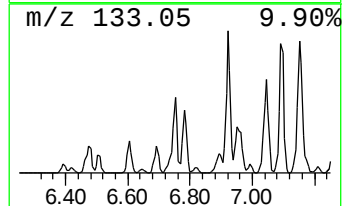
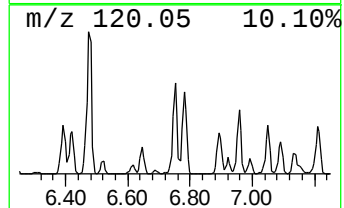
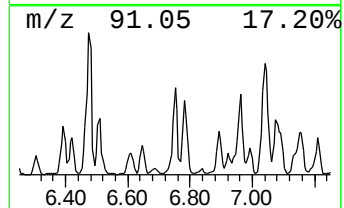
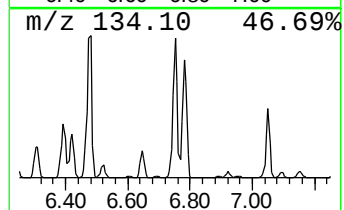
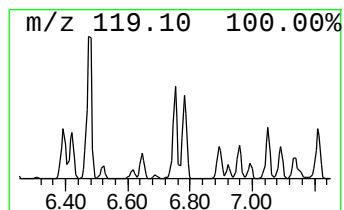
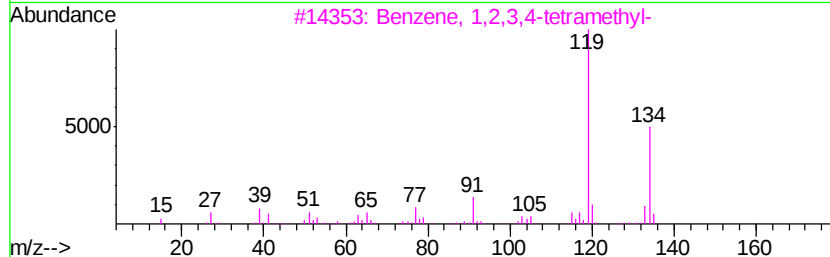
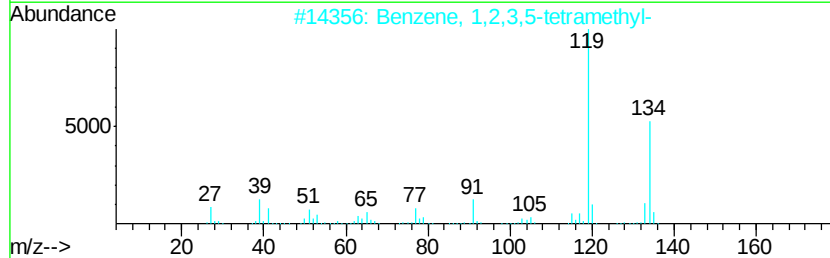
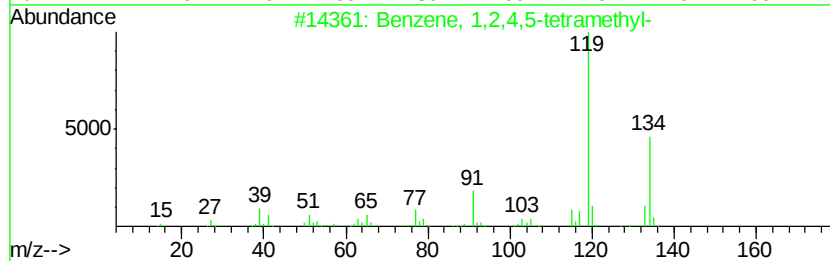
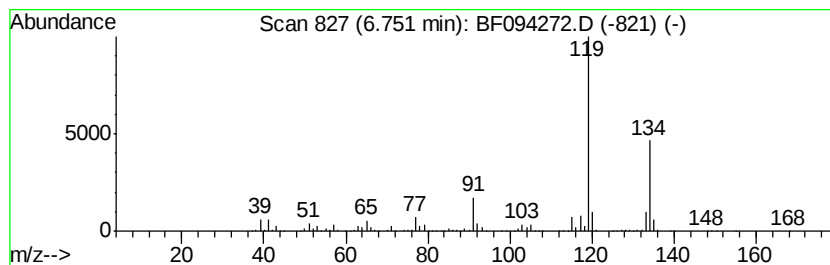
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	11.18 ng	1345520	Naphthalene-d8	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

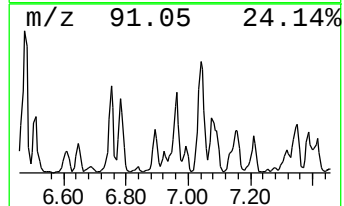
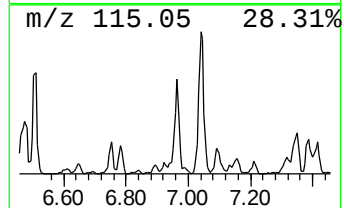
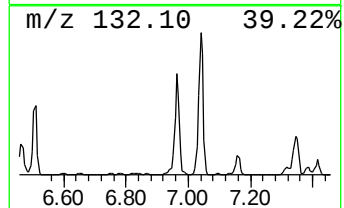
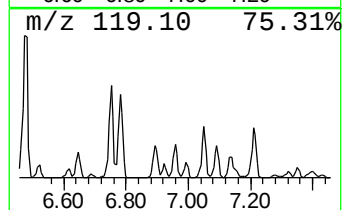
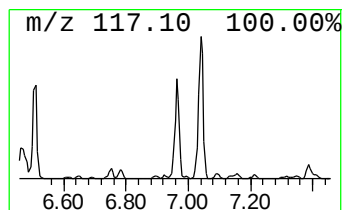
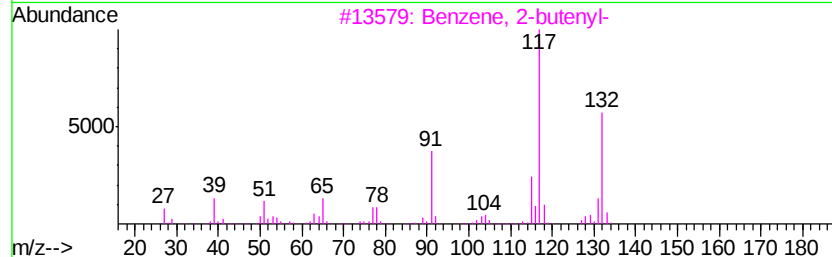
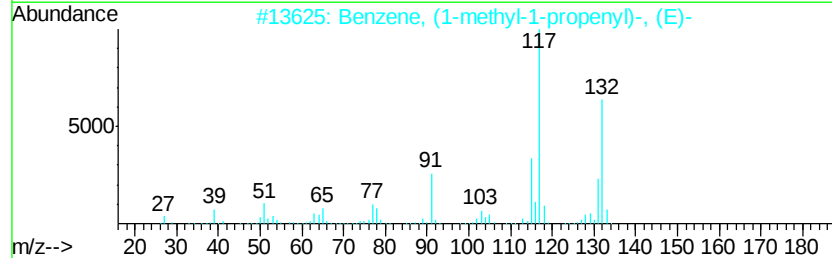
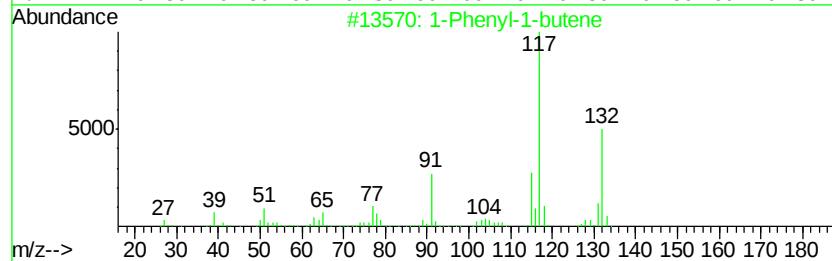
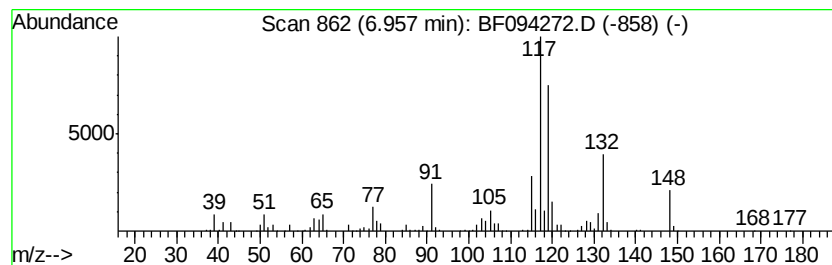
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 1-Phenyl-1-butene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.96	13.55 ng	1631480	Naphthalene-d8	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	83
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
Data File : BF094272.D
Acq On : 7 Apr 2017 14:51
Operator : SJ/MA
Sample : I2528-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TS-6

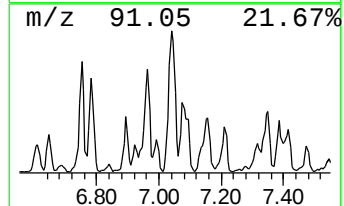
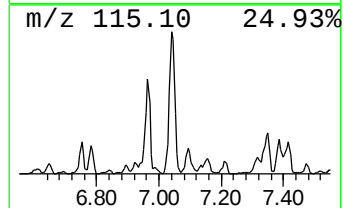
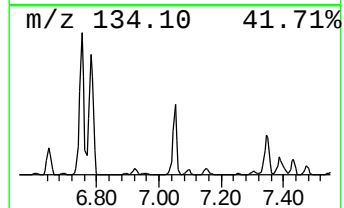
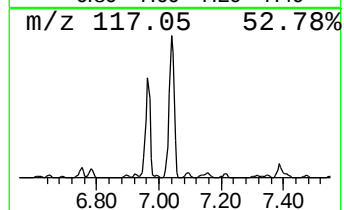
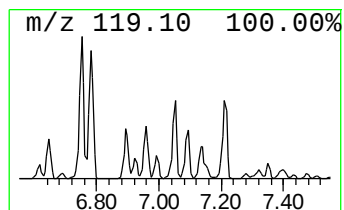
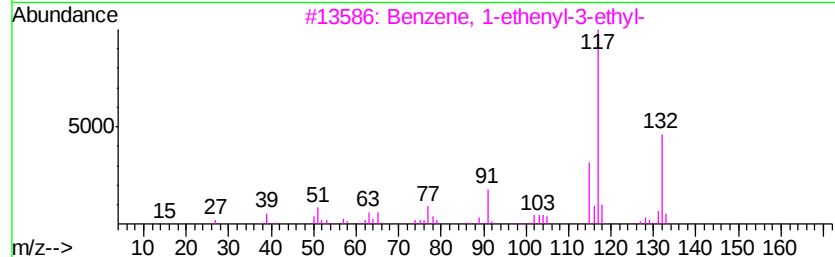
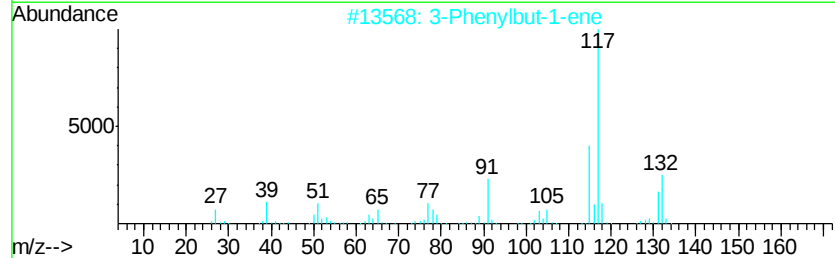
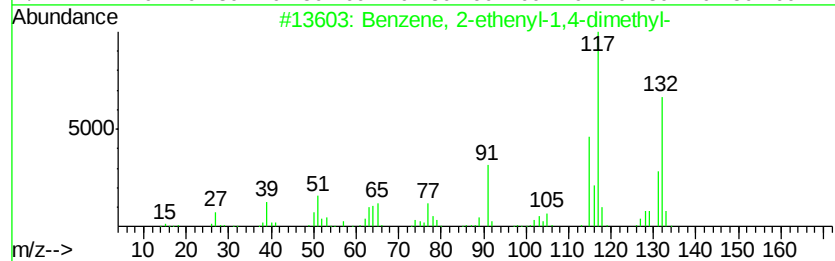
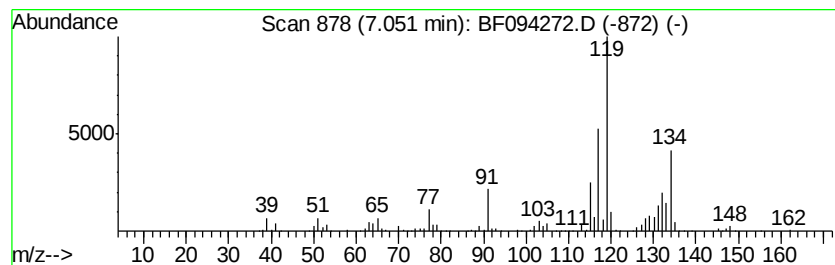
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	22.36 ng	2691840	Naphthalene-d8	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040717\
 Data File : BF094272.D
 Acq On : 7 Apr 2017 14:51
 Operator : SJ/MA
 Sample : I2528-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.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
Nonane	4.67	9.3	ng	536615	1	5.85	1158050	20.0
Benzene, propyl-	5.25	17.7	ng	1027340	1	5.85	1158050	20.0
Nonane, 2-methyl-	5.33	10.0	ng	580213	1	5.85	1158050	20.0
Benzene, 1-ethyl-...	5.35	11.6	ng	670243	1	5.85	1158050	20.0
unknown5.59	5.59	62.3	ng	3605990	1	5.85	1158050	20.0
Benzene, 1,2,3-tr...	5.65	35.4	ng	2052190	1	5.85	1158050	20.0
Benzene, 1-ethyl-...	5.92	18.0	ng	1039990	1	5.85	1158050	20.0
Indane	6.05	20.9	ng	1210470	1	5.85	1158050	20.0
Benzene, 1,2-diet...	6.14	13.5	ng	781618	1	5.85	1158050	20.0
Benzene, 1-methyl...	6.17	11.8	ng	683881	1	5.85	1158050	20.0
Benzene, 1,4-diet...	6.22	32.3	ng	1870720	1	5.85	1158050	20.0
Benzene, 1-methyl...	6.31	13.0	ng	754500	1	5.85	1158050	20.0
Benzene, 2-ethyl-...	6.39	9.7	ng	560147	1	5.85	1158050	20.0
Undecane	6.56	10.2	ng	592912	1	5.85	1158050	20.0
Benzene, 1,2,4,5-...	6.75	11.2	ng	1345520	2	7.35	2407800	20.0
1-Phenyl-1-butene	6.96	13.6	ng	1631480	2	7.35	2407800	20.0
Benzene, 2-etheny...	7.05	22.4	ng	2691840	2	7.35	2407800	20.0