

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031815\
 Data File : BF077825.D
 Acq On : 19 Mar 2015 7:07
 Operator : TP/IZ
 Sample : G1535-02 2X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP2B-6B-7B-10(0-6)

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-BF031115.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	1.298	34	36	39	rVB	1650532	2268299	100.00%	15.132%
2	1.435	45	48	57	rVB2	65356	130891	5.77%	0.873%
3	1.595	60	62	67	rVV	27605	48626	2.14%	0.324%
4	1.675	67	69	77	rVV	36539	89635	3.95%	0.598%
5	1.790	77	79	99	rVB	681809	1162630	51.26%	7.756%
6	3.264	204	208	215	rBV4	14748	49380	2.18%	0.329%
7	3.470	223	226	232	rVB	14990	36204	1.60%	0.242%
8	4.144	281	285	289	rVB2	36701	64916	2.86%	0.433%
9	4.819	341	344	348	rVB	29747	47641	2.10%	0.318%
10	4.922	352	353	357	rBV	50667	70867	3.12%	0.473%
11	5.162	371	374	376	rBV2	17908	24877	1.10%	0.166%
12	5.253	380	382	383	rVV	58067	79759	3.52%	0.532%
13	5.276	383	384	386	rVV	21784	33991	1.50%	0.227%
14	5.310	386	387	389	rVB	36757	38017	1.68%	0.254%
15	5.413	391	396	398	rBV2	112614	236720	10.44%	1.579%
16	5.459	398	400	403	rBV	285075	325672	14.36%	2.173%
17	5.619	412	414	417	rBV	18237	28206	1.24%	0.188%
18	5.710	420	422	428	rVV2	40808	74536	3.29%	0.497%
19	5.813	428	431	433	rVB	76092	84008	3.70%	0.560%
20	5.939	438	442	445	rVB	16660	27636	1.22%	0.184%
21	6.019	445	449	453	rVB5	7965	26237	1.16%	0.175%
22	6.122	453	458	460	rBV3	28253	68170	3.01%	0.455%
23	6.236	466	468	472	rBV3	43262	70567	3.11%	0.471%
24	6.305	472	474	477	rBV2	15595	22887	1.01%	0.153%
25	6.487	484	490	492	rBV2	77222	128679	5.67%	0.858%
26	6.579	492	498	499	rVV4	107598	194708	8.58%	1.299%
27	6.613	499	501	504	rVV	116727	134412	5.93%	0.897%
28	6.670	504	506	509	rVV	173232	170243	7.51%	1.136%
29	6.739	510	512	514	rVV	417189	409817	18.07%	2.734%
30	6.773	514	515	519	rVV2	93144	87979	3.88%	0.587%
31	6.876	522	524	526	rVV	457686	427602	18.85%	2.853%
32	6.945	528	530	532	rVV	431236	409876	18.07%	2.734%
33	6.979	532	533	535	rVB	92676	71452	3.15%	0.477%
34	7.105	541	544	546	rBV2	19921	46565	2.05%	0.311%

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Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	7.162	546	549	551 rVB	252208	225924	9.96%	1.507%
36	7.196	551	552	554 rBV2	33120	40932	1.80%	0.273%
37	7.230	554	555	557 rVB	126039	104422	4.60%	0.697%
38	7.345	563	565	567 rBV	344687	305620	13.47%	2.039%
39	7.379	567	568	571 rVB	90402	74354	3.28%	0.496%
40	7.482	571	577	578 rBV3	67246	89599	3.95%	0.598%
41	7.516	578	580	582 rVB2	71069	97430	4.30%	0.650%
42	7.562	582	584	586 rBV	189940	224579	9.90%	1.498%
43	7.607	586	588	589 rVB	20537	27438	1.21%	0.183%
44	7.653	589	592	596 rVB2	64700	93472	4.12%	0.624%
45	7.745	598	600	601 rBV	81112	66135	2.92%	0.441%
46	7.768	601	602	605 rVB	40444	52690	2.32%	0.352%
47	7.825	605	607	609 rBV	142238	273607	12.06%	1.825%
48	7.859	609	610	612 rVV	284761	222342	9.80%	1.483%
49	7.928	614	616	618 rBV	70079	69339	3.06%	0.463%
50	7.973	618	620	622 rVB2	30544	34035	1.50%	0.227%
51	8.008	622	623	625 rBV	26305	31317	1.38%	0.209%
52	8.122	630	633	634 rBV	118778	111480	4.91%	0.744%
53	8.156	634	636	638 rVB	122220	142518	6.28%	0.951%
54	8.270	644	646	647 rBV	33789	38676	1.71%	0.258%
55	8.293	647	648	650 rVV2	28757	43486	1.92%	0.290%
56	8.339	650	652	654 rVV	109914	112726	4.97%	0.752%
57	8.373	654	655	657 rVV2	38799	39659	1.75%	0.265%
58	8.419	657	659	661 rVV2	165264	199787	8.81%	1.333%
59	8.476	661	664	666 rVV2	56391	81863	3.61%	0.546%
60	8.533	666	669	672 rVV3	36334	70994	3.13%	0.474%
61	8.602	672	675	676 rVV	36254	42240	1.86%	0.282%
62	8.705	678	684	685 rBV2	29071	52906	2.33%	0.353%
63	8.739	685	687	688 rVV	430145	421826	18.60%	2.814%
64	8.762	688	689	692 rVV2	147644	213043	9.39%	1.421%
65	8.808	692	693	696 rVV	44151	41804	1.84%	0.279%
66	8.865	696	698	701 rVB2	23237	38860	1.71%	0.259%
67	9.093	715	718	719 rBV2	26457	32910	1.45%	0.220%
68	9.230	728	730	734 rVB	27360	32231	1.42%	0.215%
69	9.322	734	738	740 rBV3	15644	28683	1.26%	0.191%
70	9.516	753	755	758 rVB2	18755	23181	1.02%	0.155%
71	9.619	762	764	767 rVB	81213	102295	4.51%	0.682%

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72	9.745	771	775	777 rBV	55489	56231	2.48%	0.375%
73	10.088	803	805	807 rBV	477311	487739	21.50%	3.254%
74	10.911	875	877	879 rVB	374465	356438	15.71%	2.378%
75	11.882	960	962	965 rBV2	147356	210307	9.27%	1.403%
76	12.088	978	980	983 rBV2	15094	26614	1.17%	0.178%
77	12.751	1035	1038	1040 rBV	373297	395008	17.41%	2.635%
78	14.248	1167	1169	1171 rVB	55990	64002	2.82%	0.427%
79	14.328	1173	1176	1178 rBV	21407	27824	1.23%	0.186%
80	14.534	1192	1194	1196 rBV	58971	66566	2.93%	0.444%
81	14.580	1196	1198	1201 rVB3	19043	27701	1.22%	0.185%
82	14.740	1210	1212	1215 rBV	501667	616759	27.19%	4.115%
83	14.900	1224	1226	1229 rBV3	16118	37662	1.66%	0.251%
84	15.083	1239	1242	1244 rBV4	24171	43249	1.91%	0.289%
85	15.288	1257	1260	1262 rVB4	17810	31373	1.38%	0.209%
86	15.345	1262	1265	1266 rBV3	11971	26965	1.19%	0.180%
87	15.460	1273	1275	1277 rVB	20248	25032	1.10%	0.167%
88	15.517	1278	1280	1282 rVB2	20632	26053	1.15%	0.174%
89	15.940	1315	1317	1320 rBV	42723	58487	2.58%	0.390%
90	16.043	1323	1326	1331 rBV	451486	600379	26.47%	4.005%
91	16.351	1352	1353	1358 rVB3	27865	44945	1.98%	0.300%
92	16.763	1387	1389	1392 rBV3	35789	64705	2.85%	0.432%
93	17.368	1440	1442	1447 rBV	55673	133380	5.88%	0.890%
94	17.837	1480	1483	1487 rBV	300404	466214	20.55%	3.110%

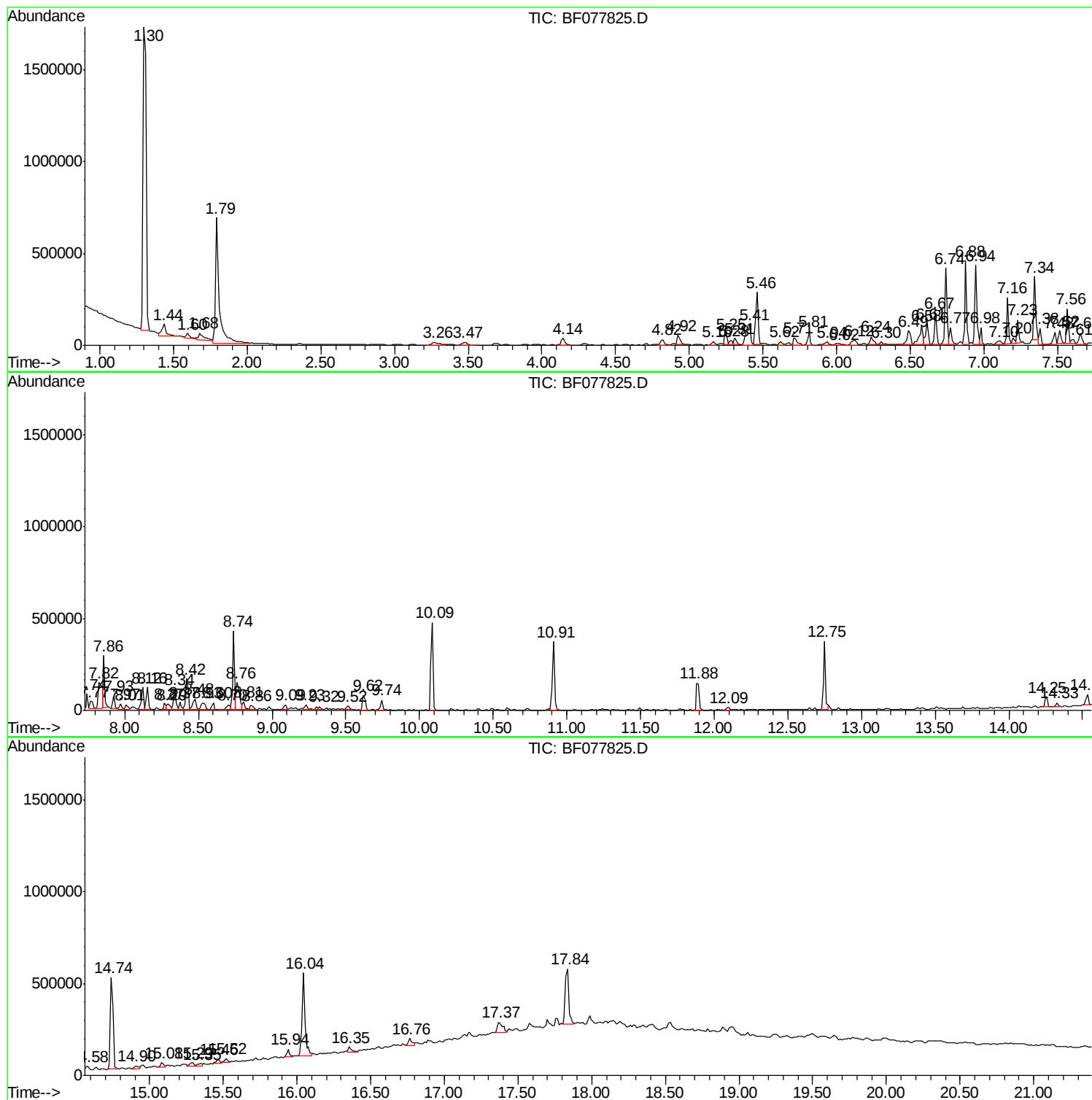
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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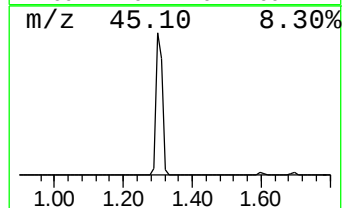
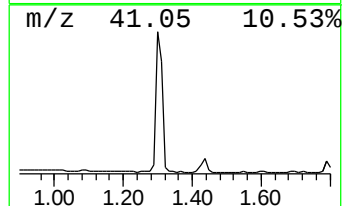
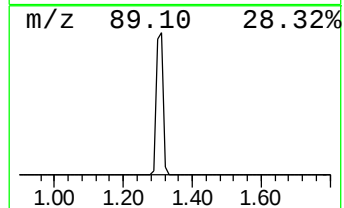
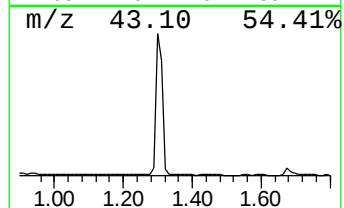
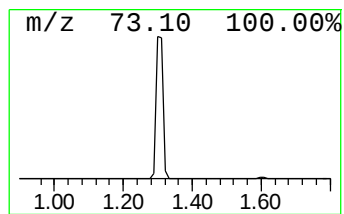
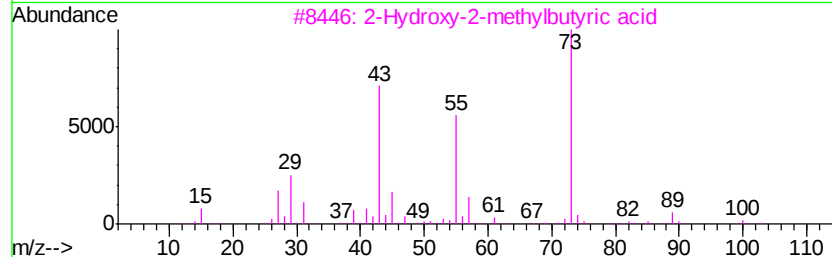
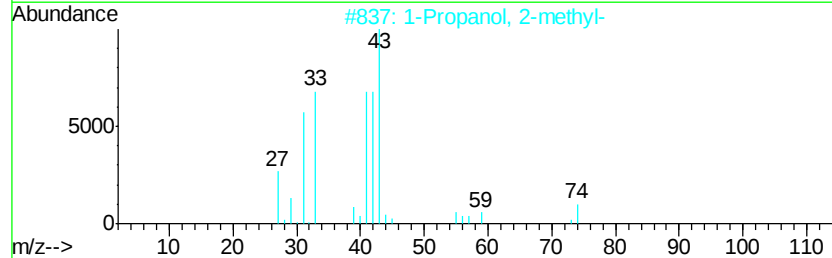
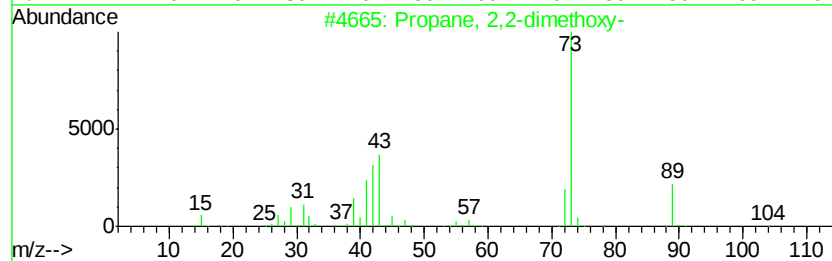
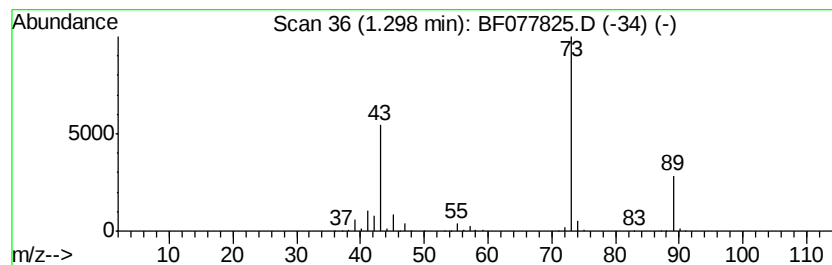
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.30	200.80 ng	2268300	1,4-Dichlorobenzene-d4	7.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9



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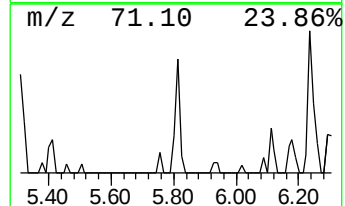
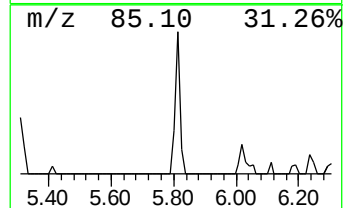
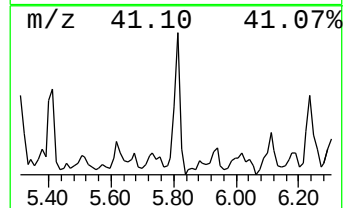
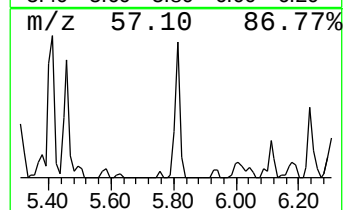
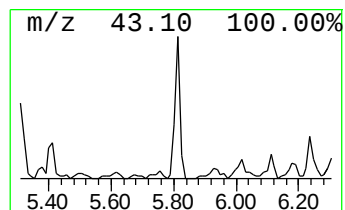
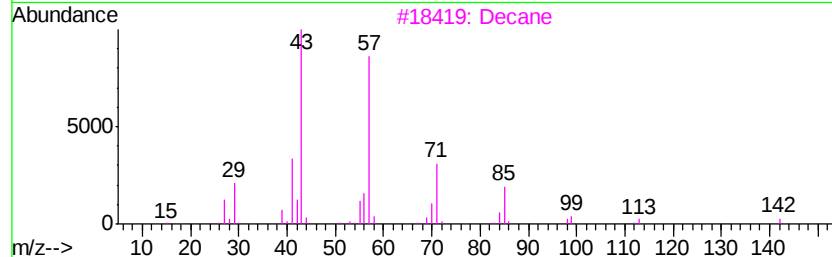
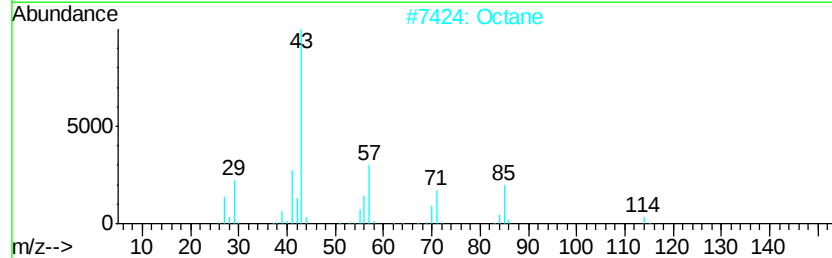
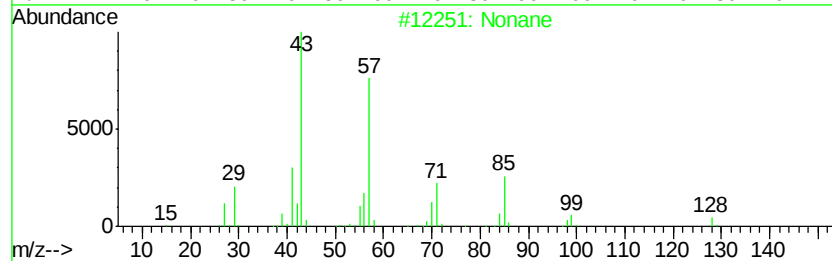
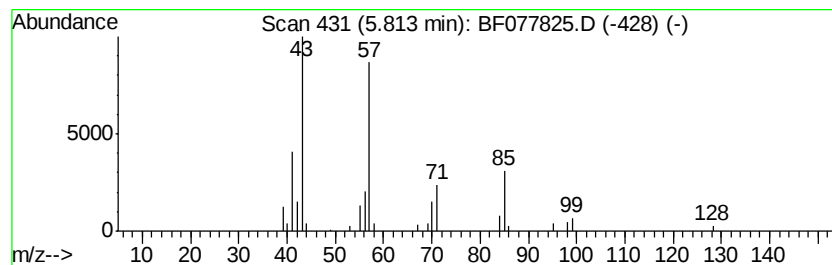
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Nonane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.81	7.44 ng	84008	1,4-Dichlorobenzene-d4	7.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonane	128	C9H20	000111-84-2	91
2		Octane	114	C8H18	000111-65-9	72
3		Decane	142	C10H22	000124-18-5	64
4		Octane, 4-methyl-	128	C9H20	002216-34-4	58
5		Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	53



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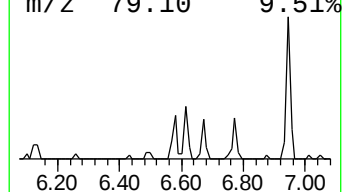
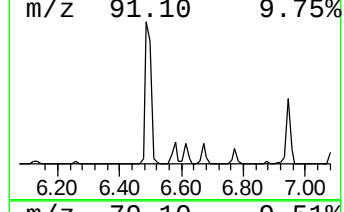
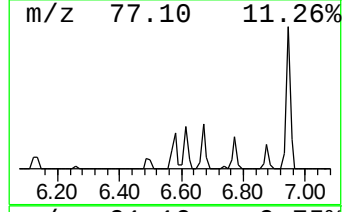
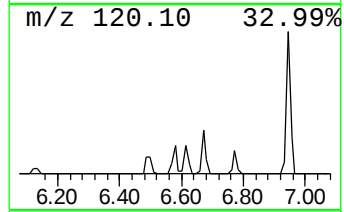
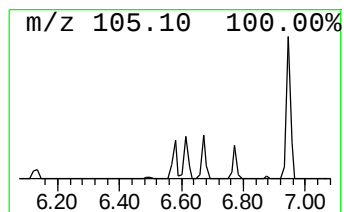
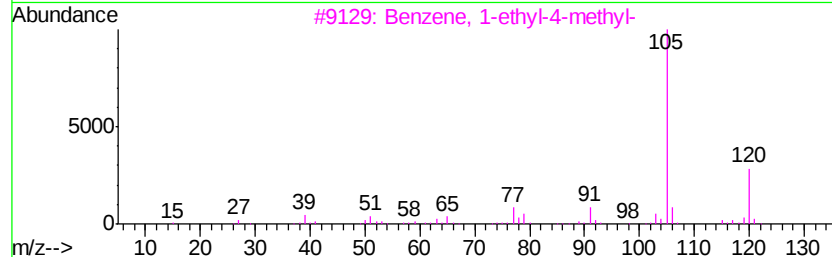
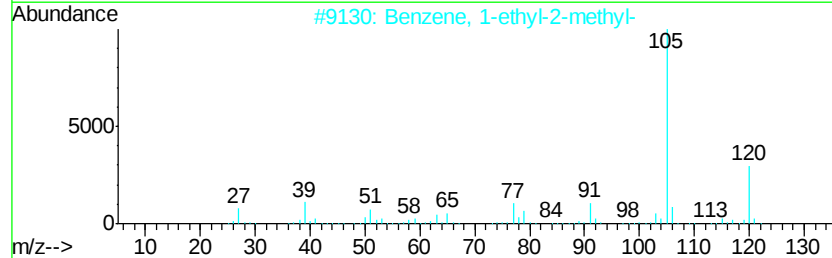
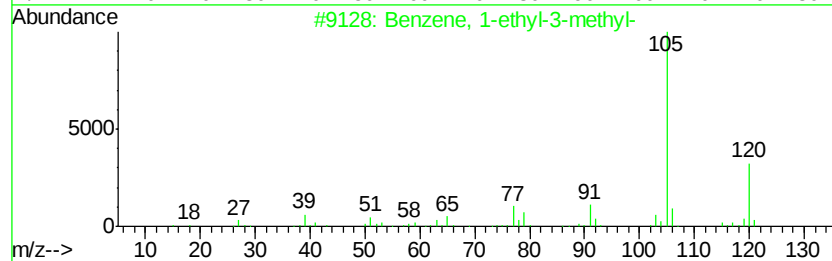
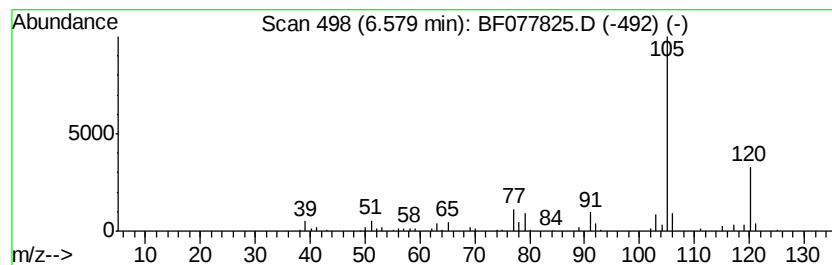
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	17.24 ng	194708	1,4-Dichlorobenzene-d4	7.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



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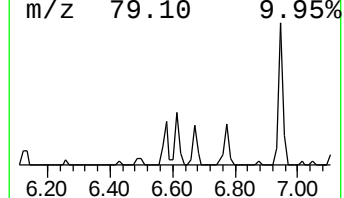
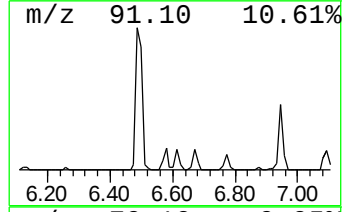
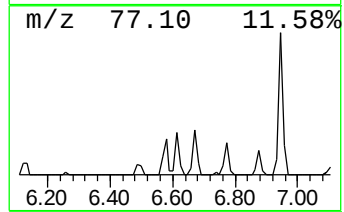
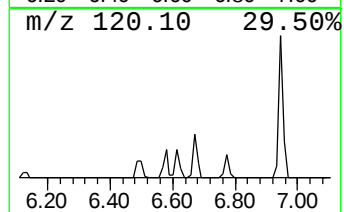
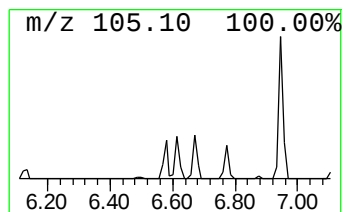
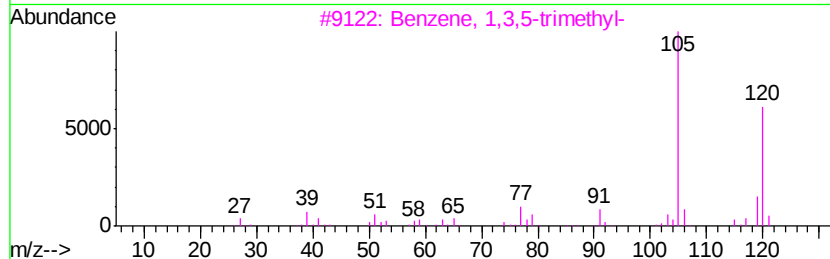
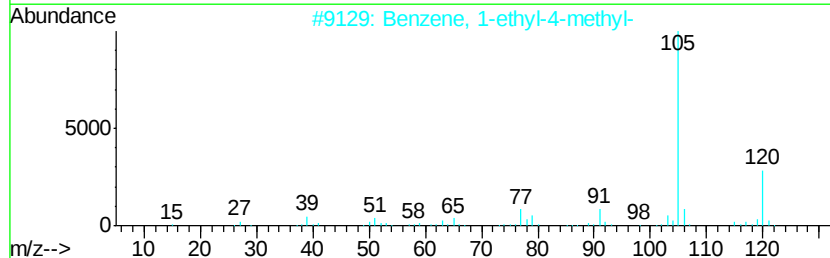
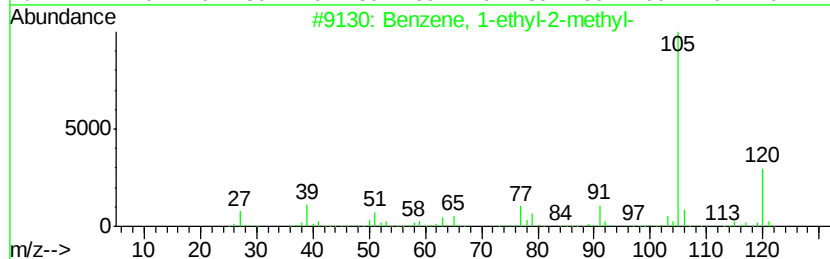
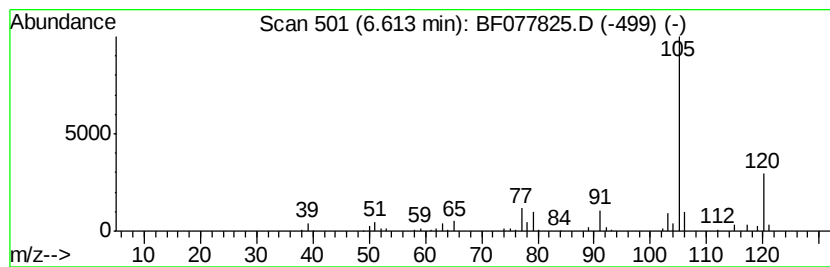
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	11.90 ng	134412	1,4-Dichlorobenzene-d4	7.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3			Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031815\
Data File : BF077825.D
Acq On : 19 Mar 2015 7:07
Operator : TP/IZ
Sample : G1535-02 2X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP2B-6B-7B-10(0-6)

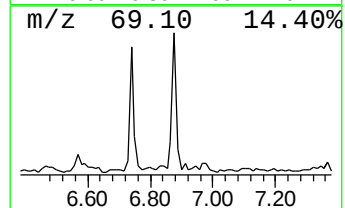
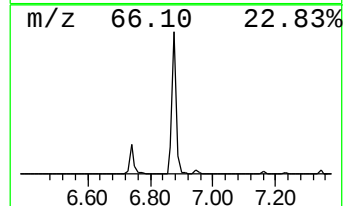
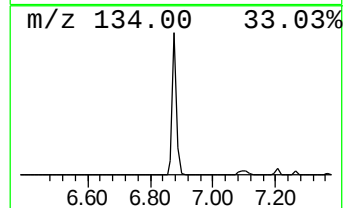
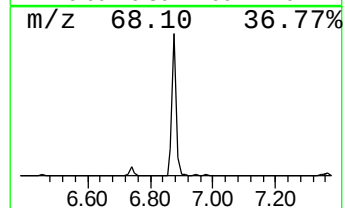
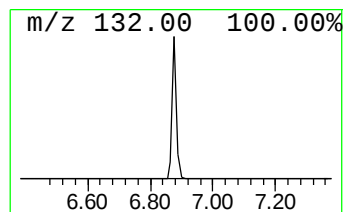
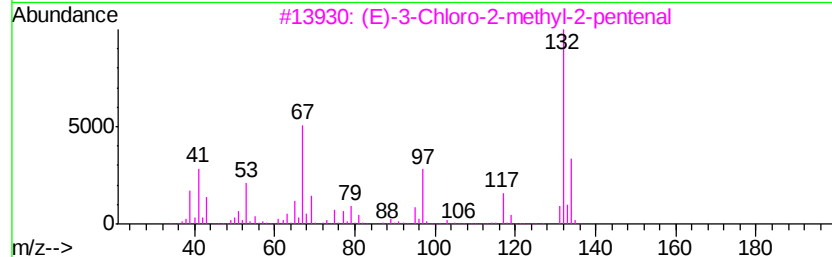
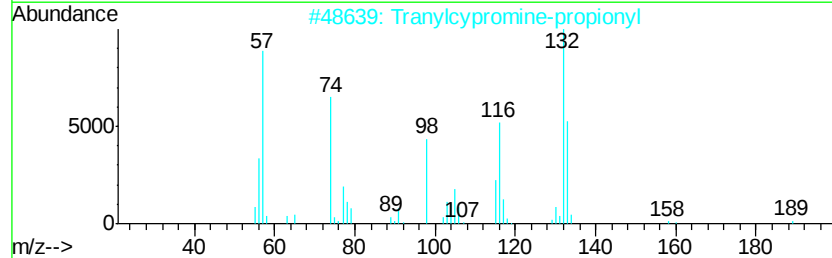
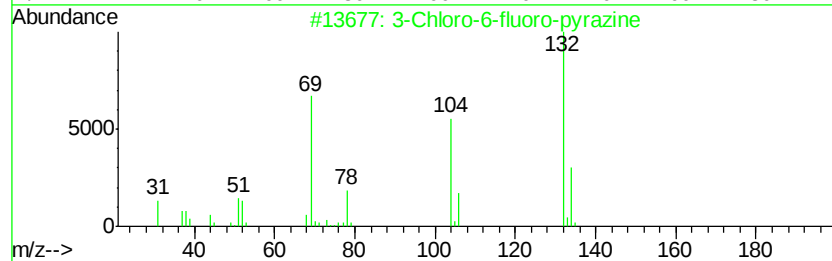
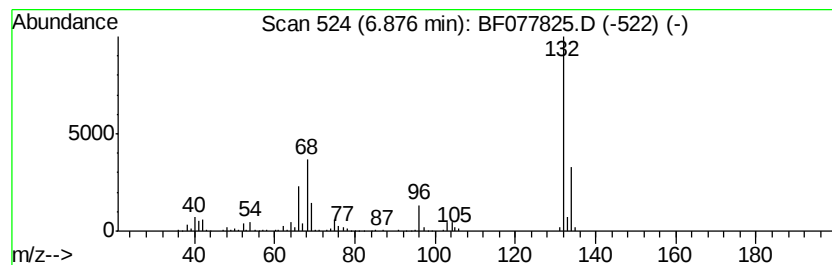
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown6.88 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	37.85 ng	427602	1,4-Dichlorobenzene-d4	7.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4			5-Aminoindole	132	C8H8N2	005192-03-0	12
5			1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031815\
Data File : BF077825.D
Acq On : 19 Mar 2015 7:07
Operator : TP/IZ
Sample : G1535-02 2X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP2B-6B-7B-10(0-6)

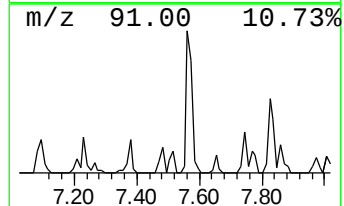
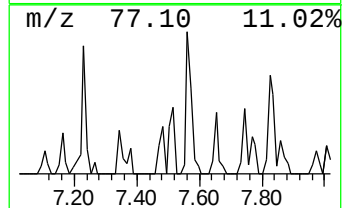
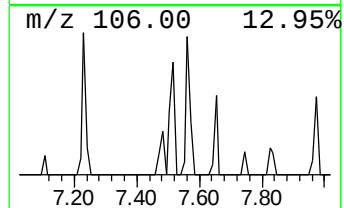
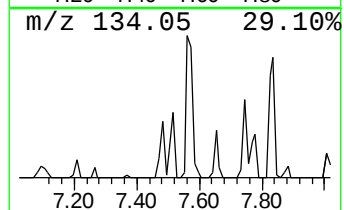
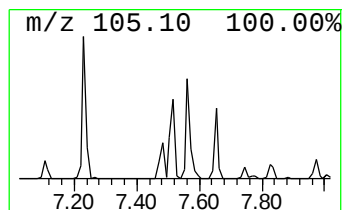
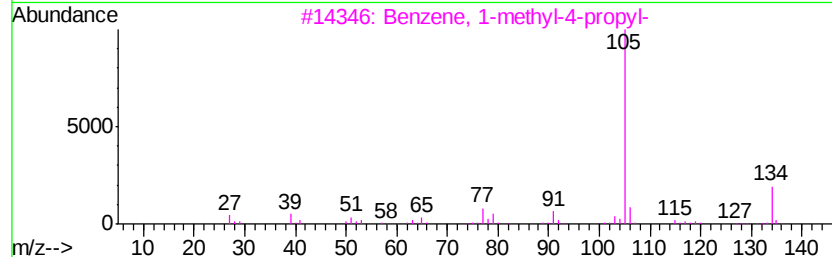
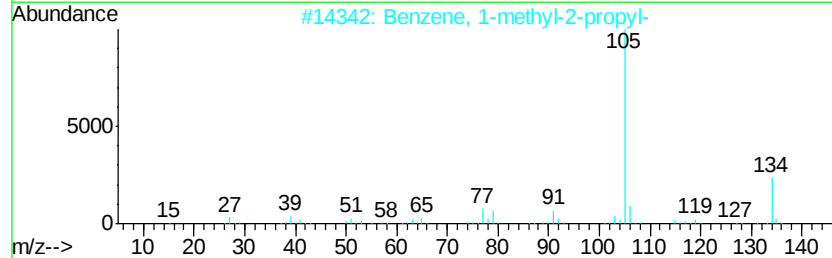
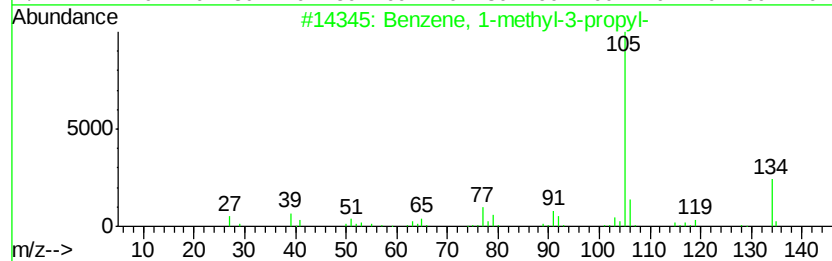
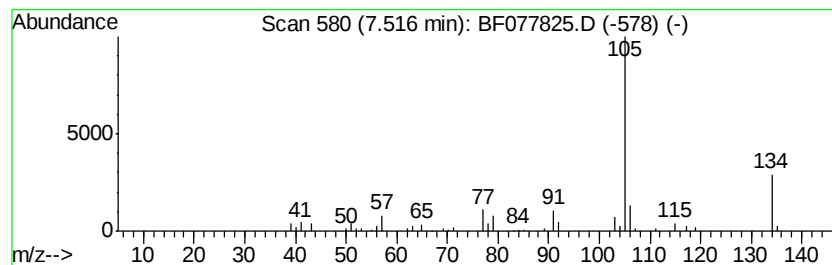
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 1-methyl-3-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	8.63 ng	97430	1,4-Dichlorobenzene-d4	7.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
2			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
5			2-Tolylloxirane	134	C9H10O	002783-26-8	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031815\
Data File : BF077825.D
Acq On : 19 Mar 2015 7:07
Operator : TP/IZ
Sample : G1535-02 2X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP2B-6B-7B-10(0-6)

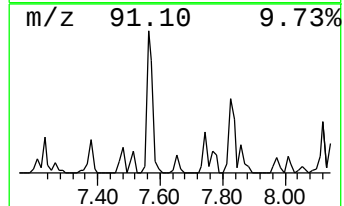
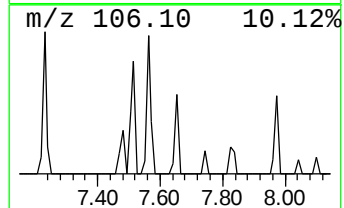
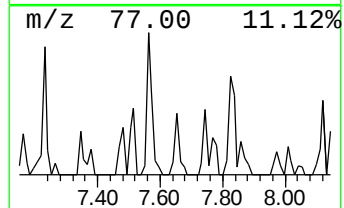
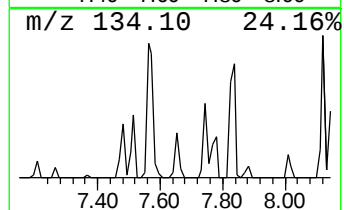
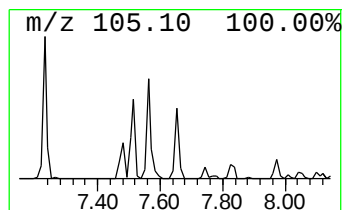
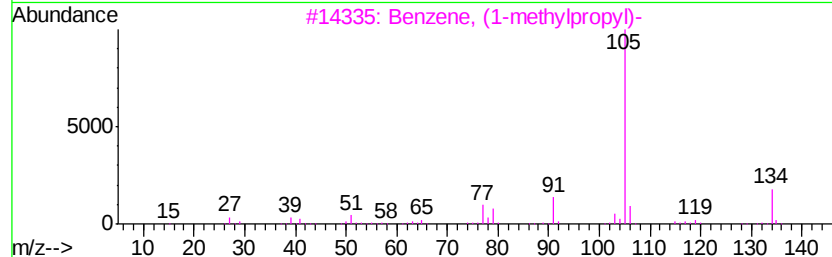
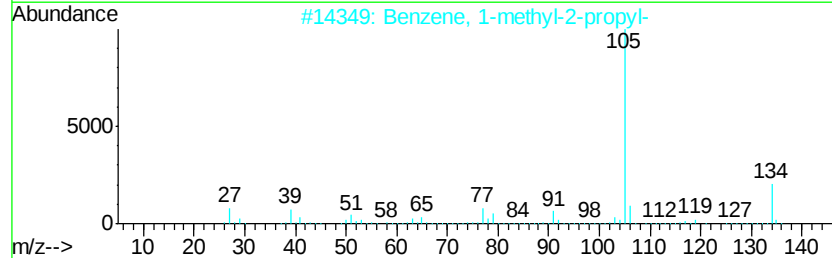
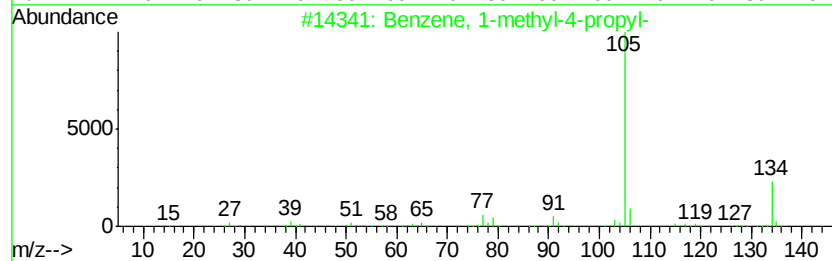
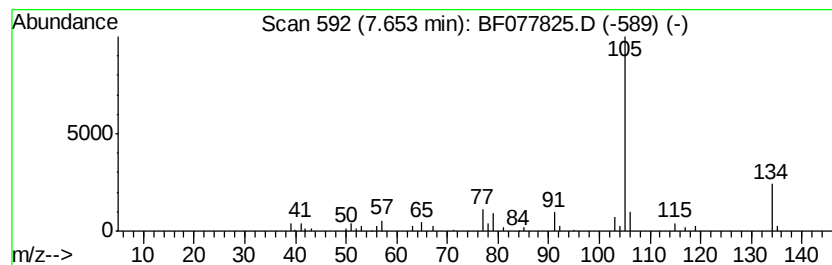
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzene, 1-methyl-4-propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	8.27 ng	93472	1,4-Dichlorobenzene-d4	7.16

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	91
3		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
5		2-Tolyloxirane	134	C9H10O	002783-26-8	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031815\
Data File : BF077825.D
Acq On : 19 Mar 2015 7:07
Operator : TP/IZ
Sample : G1535-02 2X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP2B-6B-7B-10(0-6)

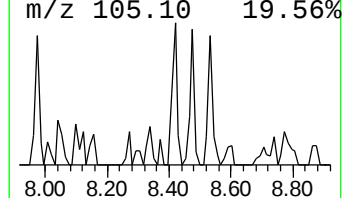
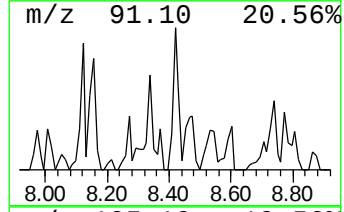
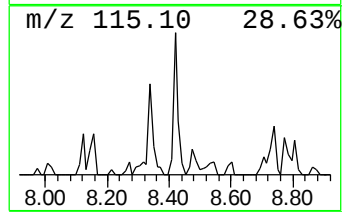
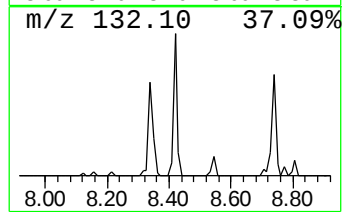
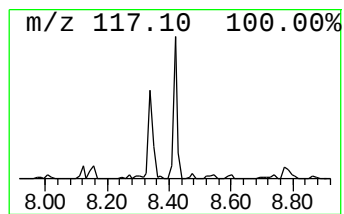
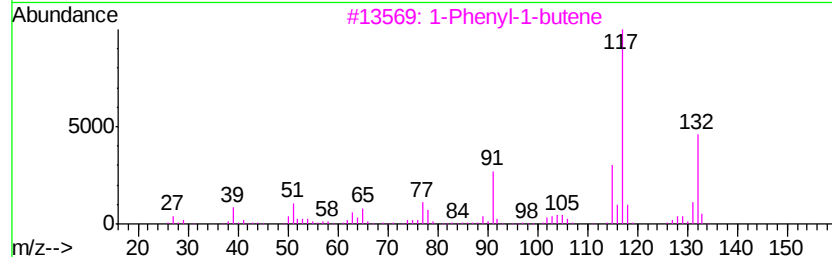
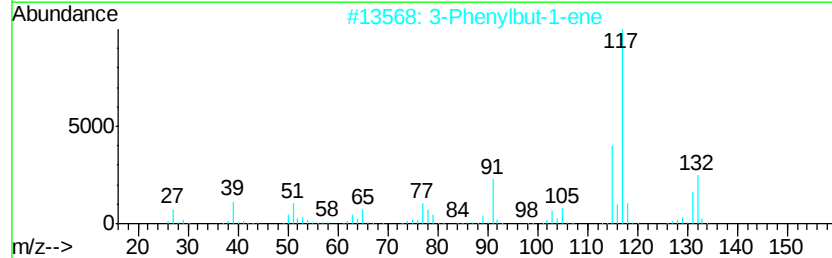
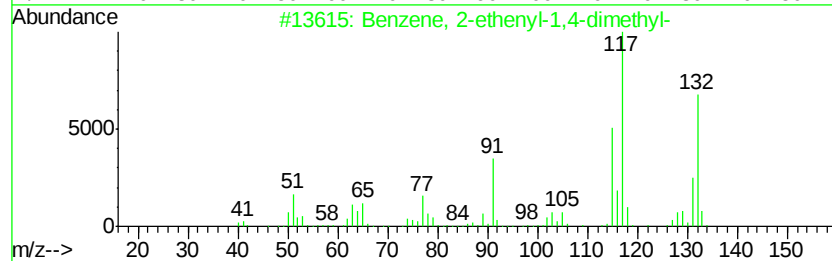
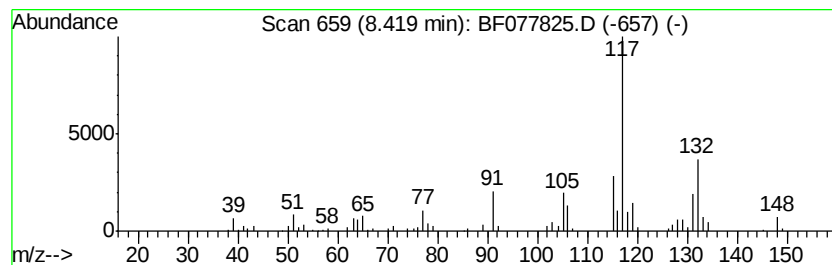
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.42	9.47 ng	199787	Napthalene-d8	8.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	76
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031815\
Data File : BF077825.D
Acq On : 19 Mar 2015 7:07
Operator : TP/IZ
Sample : G1535-02 2X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP2B-6B-7B-10(0-6)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.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
Propane, 2,2-dime...	1.30	200.8	ng	2268300	1	7.16	225924	20.0
Nonane	5.81	7.4	ng	84008	1	7.16	225924	20.0
Benzene, 1-ethyl-...	6.58	17.2	ng	194708	1	7.16	225924	20.0
Benzene, 1-ethyl-...	6.61	11.9	ng	134412	1	7.16	225924	20.0
unknown6.88	6.88	37.9	ng	427602	1	7.16	225924	20.0
Benzene, 1-methyl...	7.52	8.6	ng	97430	1	7.16	225924	20.0
Benzene, 1-methyl...	7.65	8.3	ng	93472	1	7.16	225924	20.0
Benzene, 2-etheny...	8.42	9.5	ng	199787	2	8.74	421826	20.0