

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
 Data File : BF085194.D
 Acq On : 4 Mar 2016 8:50
 Operator : SJ/UM
 Sample : H1649-02 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BT-C3(1)3W-20160301

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-BF030316.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	5.167	244	252	257	rBV	170943	219266	17.34%	1.202%
2	5.579	285	288	290	rBV	579543	738800	58.43%	4.050%
3	5.819	307	309	311	rBV	57801	47104	3.73%	0.258%
4	6.139	333	337	342	rBV5	12265	34949	2.76%	0.192%
5	6.230	342	345	348	rBV2	13003	24416	1.93%	0.134%
6	6.436	358	363	364	rBV3	16898	28516	2.26%	0.156%
7	6.470	364	366	368	rVV	74676	84884	6.71%	0.465%
8	6.504	368	369	370	rVV	139734	105617	8.35%	0.579%
9	6.539	370	372	373	rVV	44735	55555	4.39%	0.305%
10	6.596	373	377	379	rVV	654598	900899	71.25%	4.938%
11	6.664	381	383	384	rVV	80840	80645	6.38%	0.442%
12	6.744	384	390	392	rVV	759782	991956	78.45%	5.437%
13	6.802	393	395	398	rVB	258804	240180	19.00%	1.316%
14	6.973	408	410	413	rBV2	639343	894323	70.73%	4.902%
15	7.042	413	416	417	rVB2	43732	46248	3.66%	0.253%
16	7.133	422	424	426	rVB	778655	630780	49.89%	3.457%
17	7.224	431	432	433	rBV	17059	23190	1.83%	0.127%
18	7.259	433	435	437	rVV	82696	74188	5.87%	0.407%
19	7.293	437	438	442	rVB2	54473	107247	8.48%	0.588%
20	7.373	442	445	447	rVB2	27423	34259	2.71%	0.188%
21	7.442	447	451	452	rBV3	34300	68275	5.40%	0.374%
22	7.533	457	459	462	rVV	574930	531354	42.02%	2.912%
23	7.579	462	463	465	rVB	26099	19836	1.57%	0.109%
24	7.670	469	471	473	rBV2	23639	29910	2.37%	0.164%
25	7.750	475	478	479	rBV2	36998	41431	3.28%	0.227%
26	7.773	479	480	483	rVB	48576	44441	3.51%	0.244%
27	7.887	487	490	491	rVB2	13572	18552	1.47%	0.102%
28	7.933	491	494	497	rBV2	29696	57259	4.53%	0.314%
29	8.002	497	500	502	rBV	68418	82539	6.53%	0.452%
30	8.116	508	510	515	rBV2	15492	35077	2.77%	0.192%
31	8.265	520	523	525	rBV	1458207	1264407	100.00%	6.931%
32	8.653	555	557	561	rVB4	9286	18847	1.49%	0.103%
33	8.767	565	567	573	rBV4	9166	24152	1.91%	0.132%
34	8.973	583	585	588	rVB	38247	37265	2.95%	0.204%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.076	590	594	598 rBV	34418	43051	3.40%	0.236%
36	9.339	614	617	619 rBV	1071153	897450	70.98%	4.919%
37	9.682	642	647	650 rBV5	21770	56310	4.45%	0.309%
38	9.728	650	651	655 rVB	148664	138441	10.95%	0.759%
39	9.876	663	664	668 rBV2	31766	47922	3.79%	0.263%
40	9.945	668	670	674 rVV	40582	50936	4.03%	0.279%
41	10.025	674	677	678 rVV	869603	1176567	93.05%	6.449%
42	10.048	678	679	684 rVV3	34304	59856	4.73%	0.328%
43	10.219	692	694	696 rVV2	15957	21832	1.73%	0.120%
44	10.265	696	698	700 rVB2	21845	32051	2.53%	0.176%
45	10.448	711	714	716 rVB	57067	47827	3.78%	0.262%
46	10.665	731	733	737 rBV2	24079	59629	4.72%	0.327%
47	10.802	743	745	747 rBV	192959	178377	14.11%	0.978%
48	10.848	747	749	751 rVB3	21424	33632	2.66%	0.184%
49	10.916	754	755	761 rBV	50436	98924	7.82%	0.542%
50	11.008	761	763	765 rBV3	20571	40527	3.21%	0.222%
51	11.111	770	772	776 rVB2	30349	69051	5.46%	0.378%
52	11.362	792	794	796 rVV2	28002	47424	3.75%	0.260%
53	11.408	796	798	803 rVB2	52554	108245	8.56%	0.593%
54	11.499	804	806	810 rBV2	834886	1172711	92.75%	6.428%
55	11.579	810	813	814 rVB	52118	53514	4.23%	0.293%
56	11.728	823	826	830 rBV2	30080	45756	3.62%	0.251%
57	11.842	834	836	839 rVB4	9991	19481	1.54%	0.107%
58	11.911	840	842	845 rVB3	23193	46006	3.64%	0.252%
59	12.002	845	850	851 rBV	43489	81879	6.48%	0.449%
60	12.036	851	853	856 rVV	140688	156150	12.35%	0.856%
61	12.082	856	857	858 rVV	25540	25747	2.04%	0.141%
62	12.116	858	860	864 rVB2	61565	92120	7.29%	0.505%
63	12.322	872	878	879 rBV3	31684	65473	5.18%	0.359%
64	12.391	882	884	886 rVV	77664	114415	9.05%	0.627%
65	12.436	886	888	897 rVB2	160290	314317	24.86%	1.723%
66	12.551	897	898	900 rVB	24965	23792	1.88%	0.130%
67	12.631	904	905	909 rVB	13801	20900	1.65%	0.115%
68	12.711	910	912	914 rVB	270867	263510	20.84%	1.444%
69	12.779	914	918	919 rBV3	13177	24600	1.95%	0.135%
70	12.871	922	926	927 rBV2	43729	82642	6.54%	0.453%
71	12.905	927	929	930 rBV	153576	133125	10.53%	0.730%

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72	12.939	930	932	935	rVB	199371	238641	18.87%	1.308%
73	13.088	941	945	947	rBV	450644	585983	46.34%	3.212%
74	13.156	949	951	953	rVB2	37942	61884	4.89%	0.339%
75	13.202	953	955	956	rBV2	27466	47751	3.78%	0.262%
76	13.385	968	971	977	rBV2	38867	126359	9.99%	0.693%
77	13.476	977	979	980	rBV	27230	35433	2.80%	0.194%
78	13.568	982	987	989	rBV5	42595	131979	10.44%	0.723%
79	13.614	989	991	992	rVB	67932	70679	5.59%	0.387%
80	13.659	992	995	999	rBV5	39805	136666	10.81%	0.749%
81	13.774	1003	1005	1007	rBV	56695	79547	6.29%	0.436%
82	13.865	1011	1013	1016	rVB3	20603	37238	2.95%	0.204%
83	13.979	1021	1023	1027	rVB2	59839	84388	6.67%	0.463%
84	14.139	1034	1037	1038	rBV	985322	1003721	79.38%	5.502%
85	14.242	1044	1046	1047	rBV2	43596	37497	2.97%	0.206%
86	14.608	1075	1078	1080	rBV2	36952	79772	6.31%	0.437%
87	14.757	1089	1091	1095	rBV	44835	88443	6.99%	0.485%
88	14.997	1109	1112	1113	rBV2	35034	51140	4.04%	0.280%
89	15.088	1118	1120	1122	rBV2	36290	67638	5.35%	0.371%
90	15.179	1125	1128	1132	rVV	118474	214195	16.94%	1.174%
91	15.488	1152	1155	1157	rVB	99915	168891	13.36%	0.926%
92	15.545	1158	1160	1163	rVV	84132	114103	9.02%	0.625%
93	15.614	1163	1166	1172	rVB	655626	911039	72.05%	4.994%
94	15.945	1193	1195	1196	rBV2	40891	37742	2.98%	0.207%
95	16.837	1272	1273	1274	rVB	27080	18563	1.47%	0.102%
96	17.077	1292	1294	1299	rVB2	45058	83948	6.64%	0.460%
97	17.523	1331	1333	1337	rVB	29537	57645	4.56%	0.316%
98	18.265	1396	1398	1400	rBV2	24793	19713	1.56%	0.108%
99	19.283	1485	1487	1489	rVB2	31251	26272	2.08%	0.144%
100	19.420	1498	1499	1501	rBV2	24728	46505	3.68%	0.255%

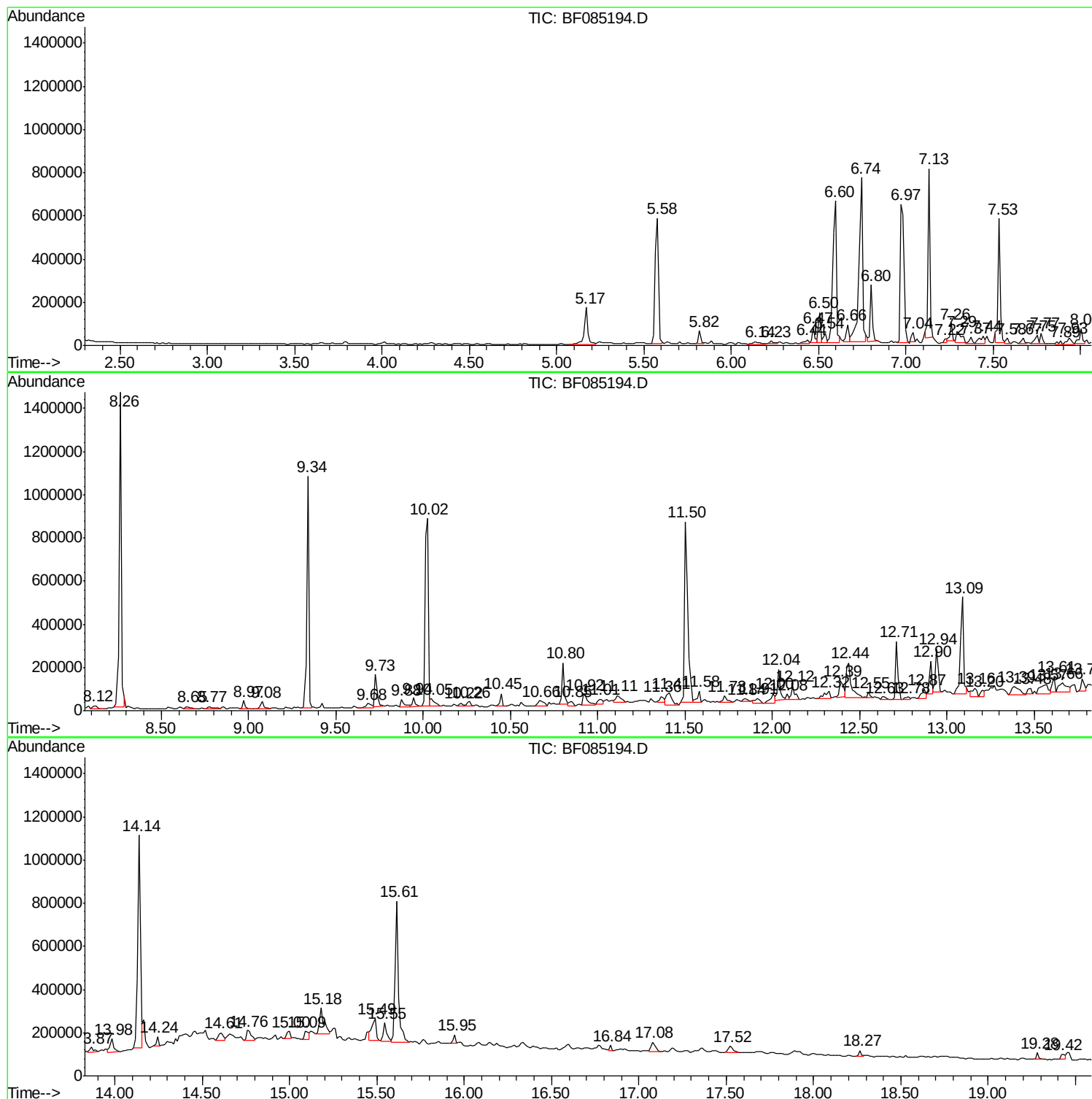
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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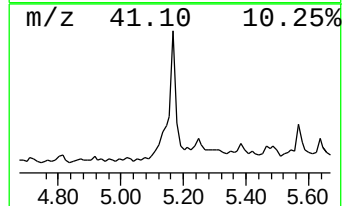
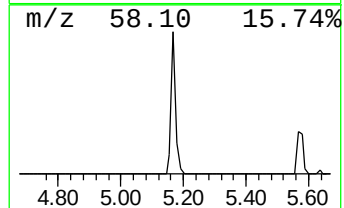
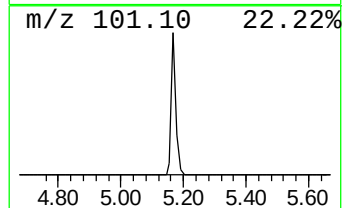
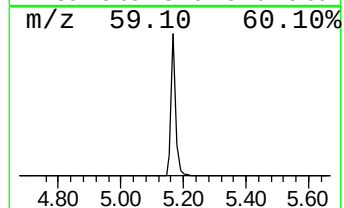
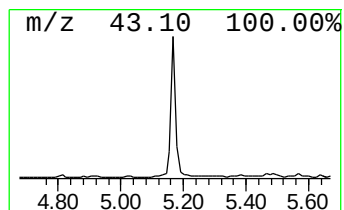
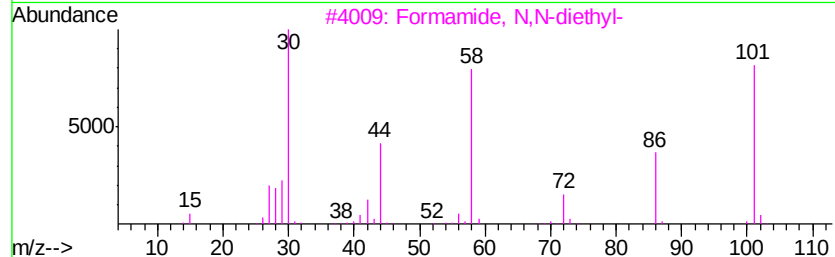
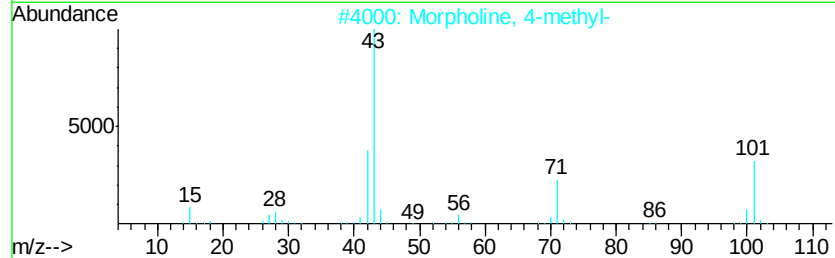
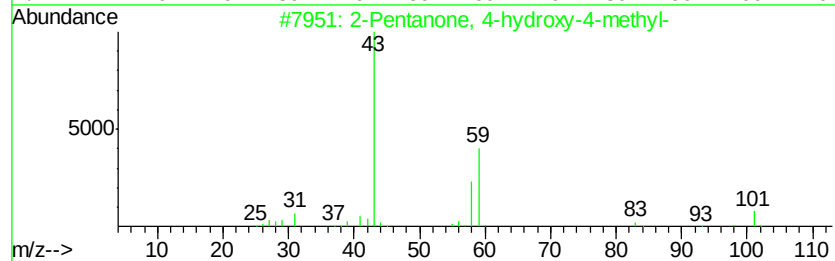
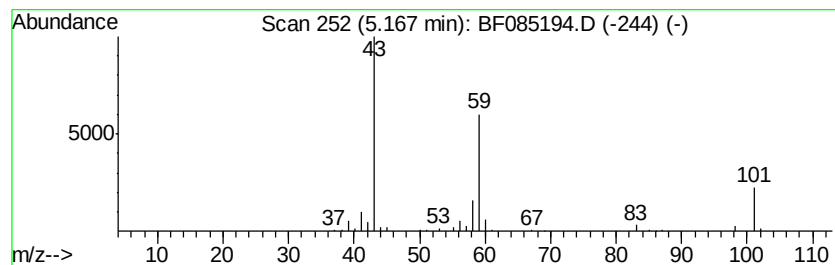
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.17	4.90 ng	219266	1,4-Dichlorobenzene-d4	6.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9



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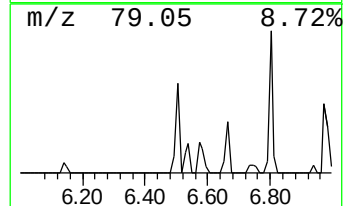
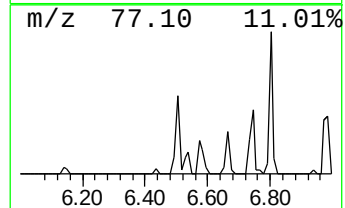
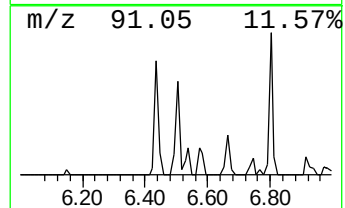
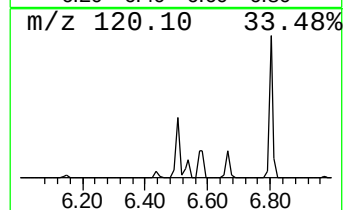
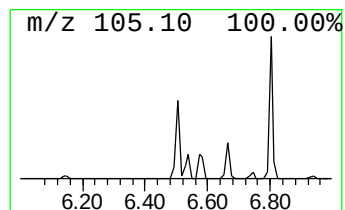
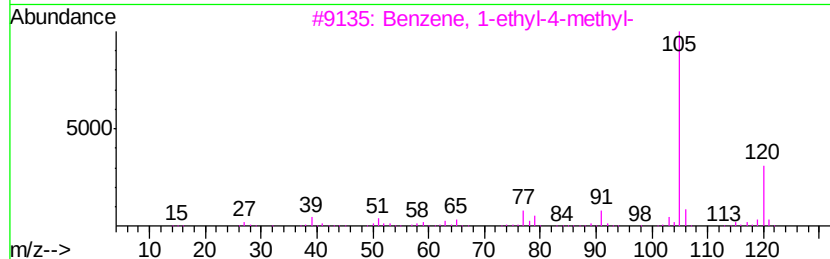
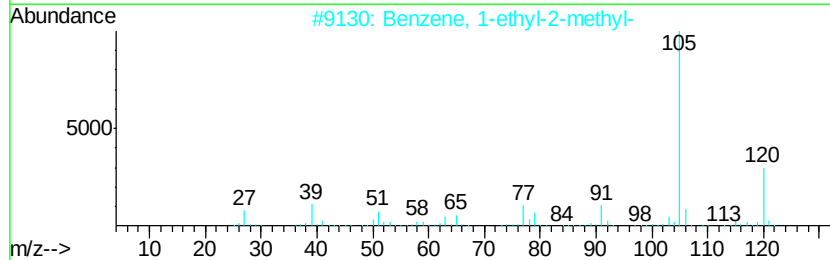
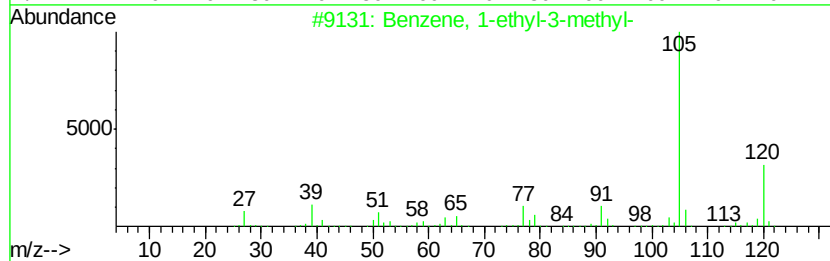
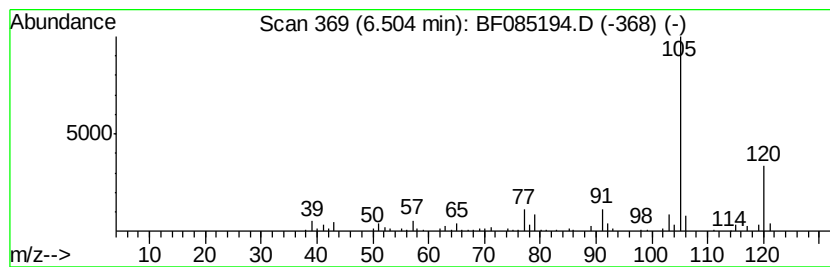
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	2.36 ng	105617	1,4-Dichlorobenzene-d4	6.98

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



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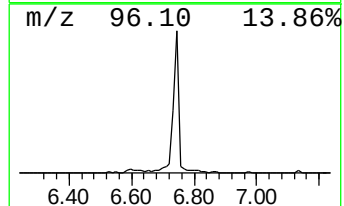
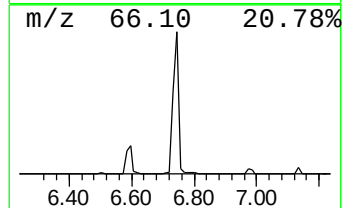
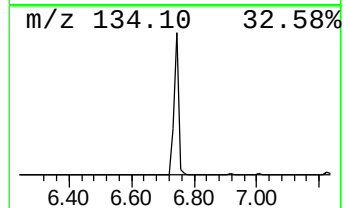
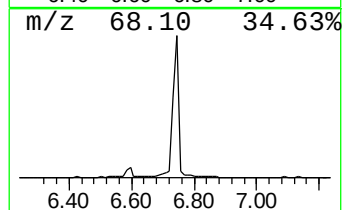
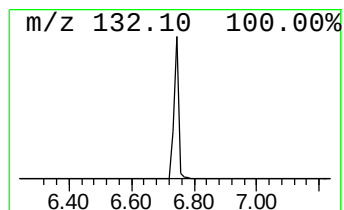
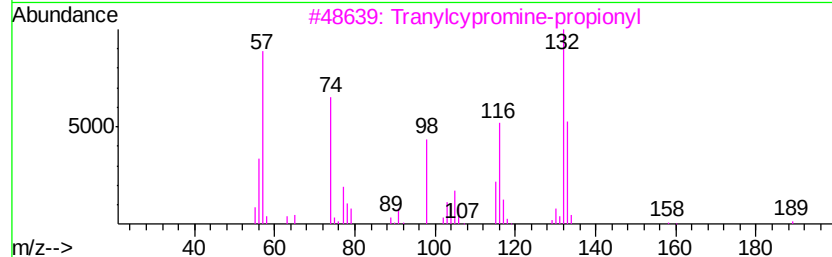
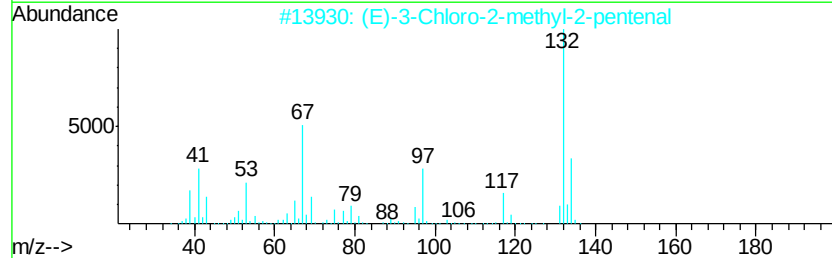
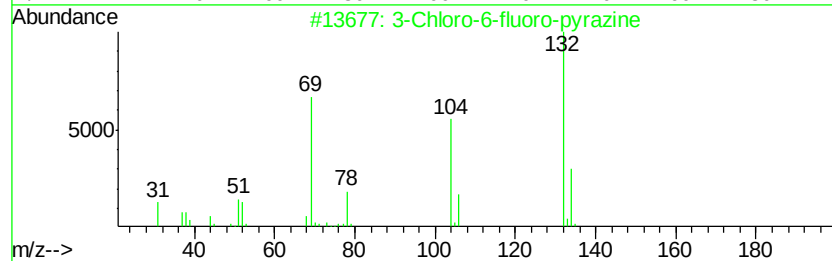
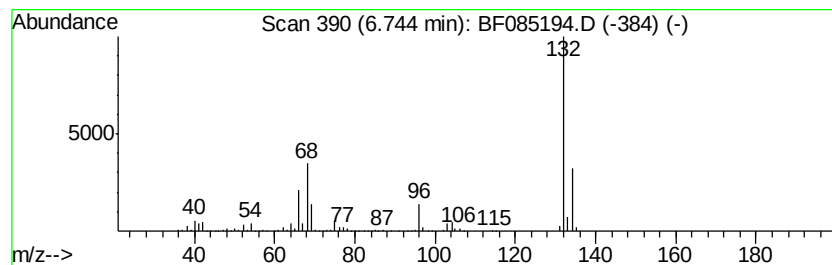
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	22.18 ng	991956	1,4-Dichlorobenzene-d4	6.98

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		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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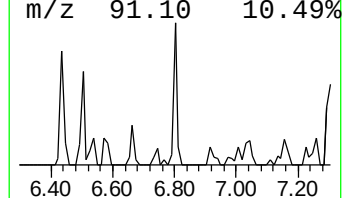
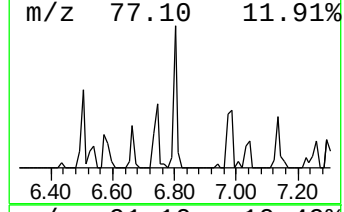
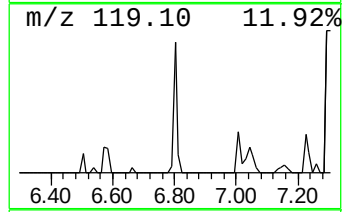
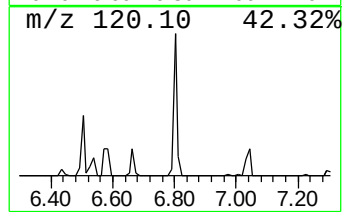
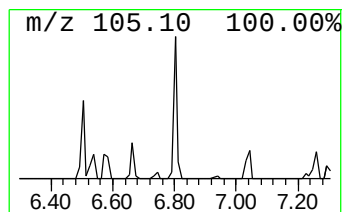
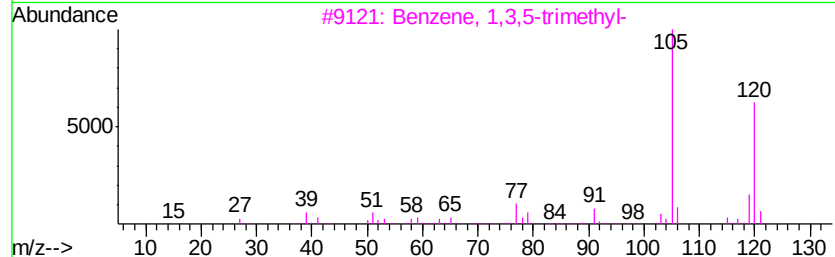
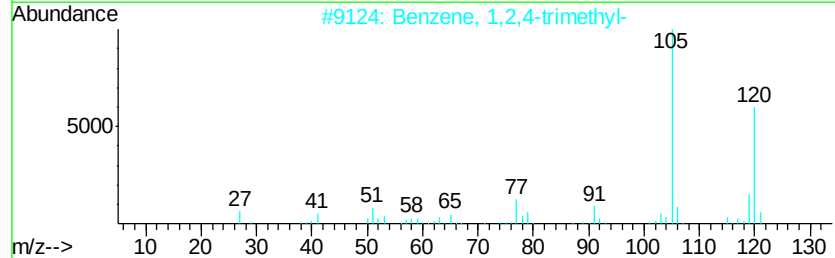
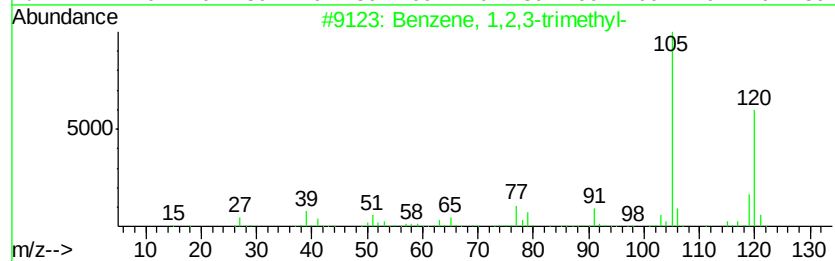
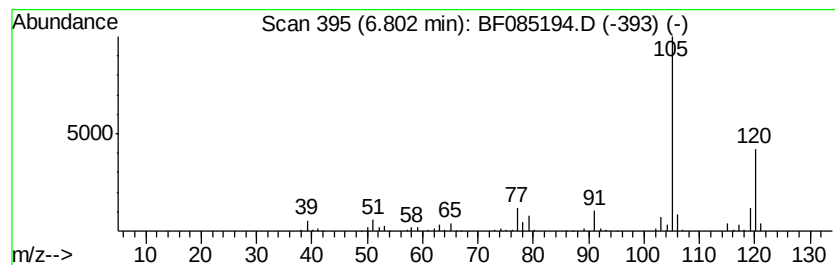
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	5.37 ng	240180	1,4-Dichlorobenzene-d4	6.98

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	95
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
Data File : BF085194.D
Acq On : 4 Mar 2016 8:50
Operator : SJ/UM
Sample : H1649-02 5X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampled :
BT-C3(1)3W-20160301

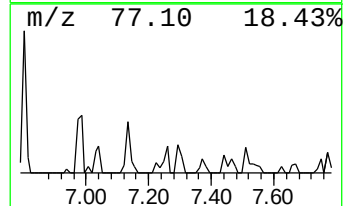
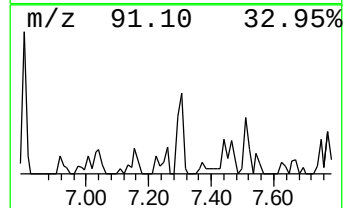
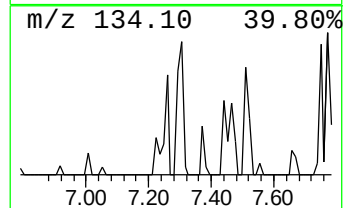
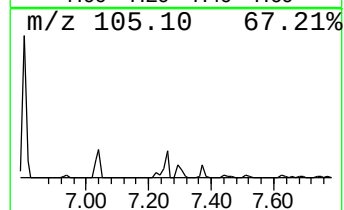
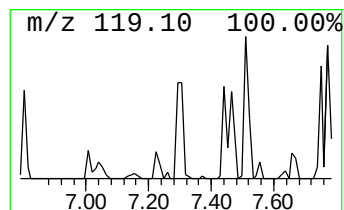
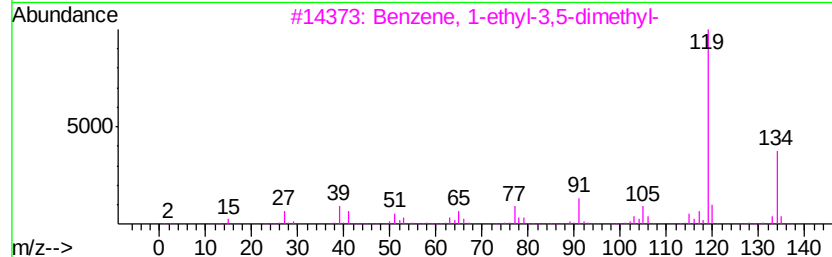
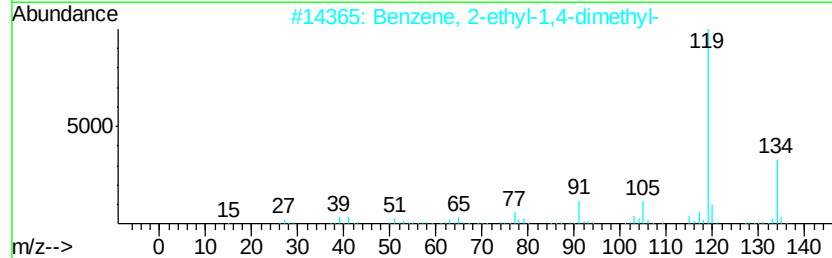
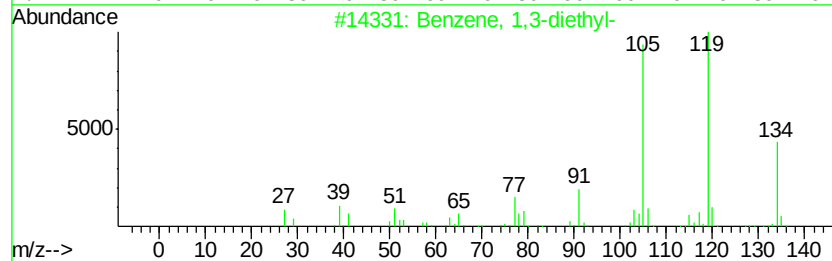
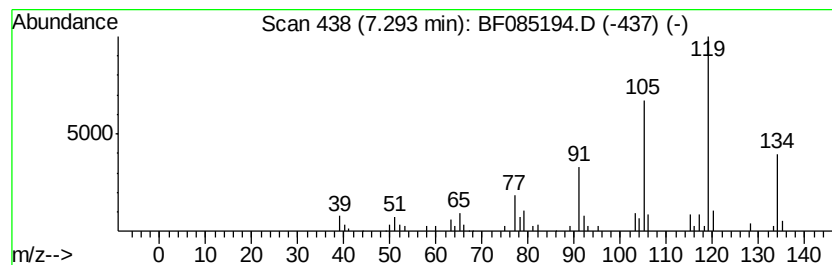
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1,3-diethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.29	2.40 ng	107247	1,4-Dichlorobenzene-d4	6.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
Data File : BF085194.D
Acq On : 4 Mar 2016 8:50
Operator : SJ/UM
Sample : H1649-02 5X
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Instrument :
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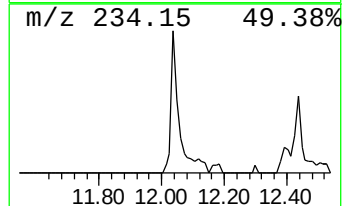
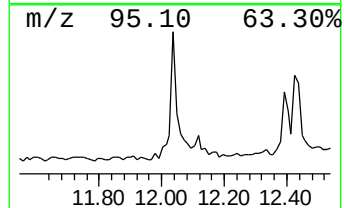
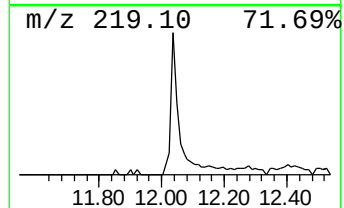
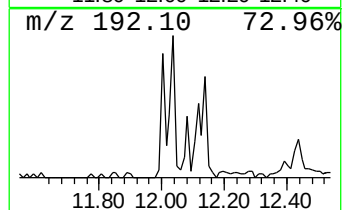
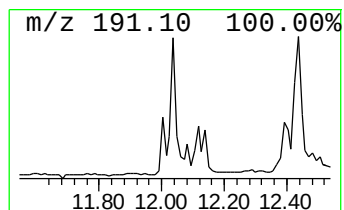
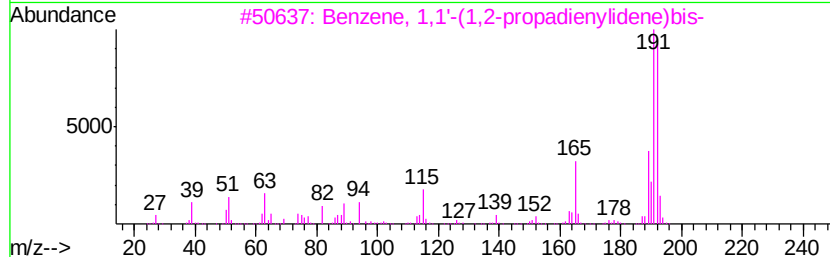
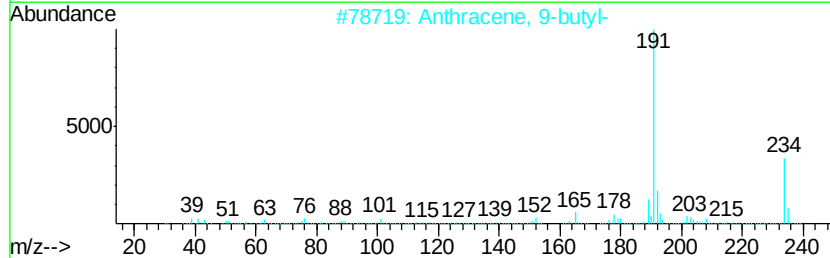
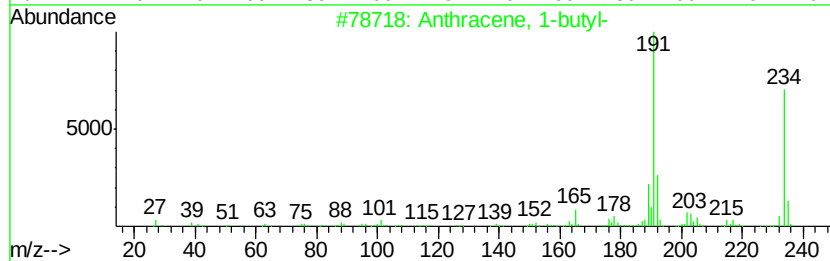
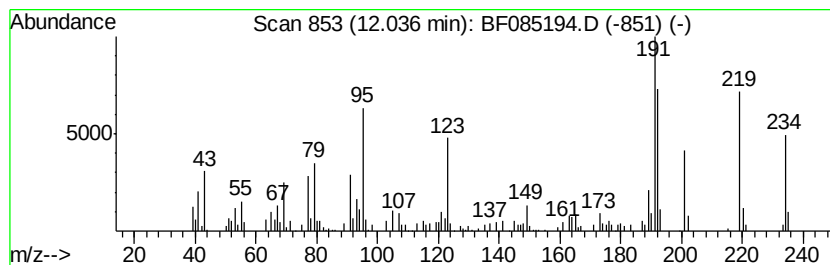
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown12.04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	2.66 ng	156150	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-butyl-	234	C18H18	1000156-51-7	46
2		Anthracene, 9-butyl-	234	C18H18	001498-69-7	38
3		Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	35
4		Phenanthrene, 9-butyl-	234	C18H18	010394-57-7	35
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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Sample : H1649-02 5X
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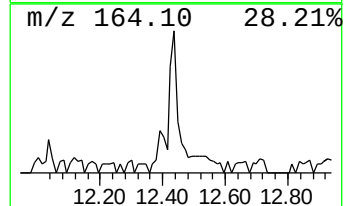
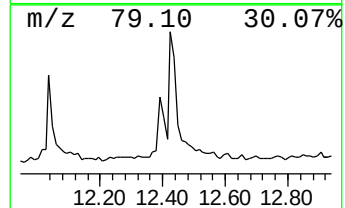
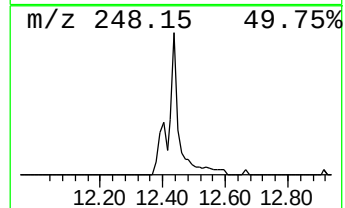
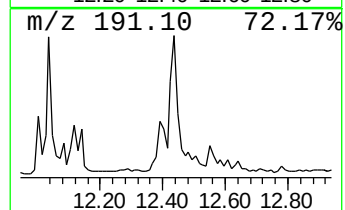
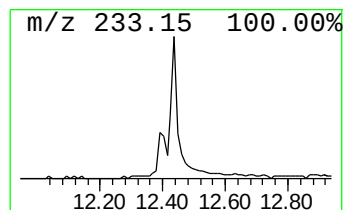
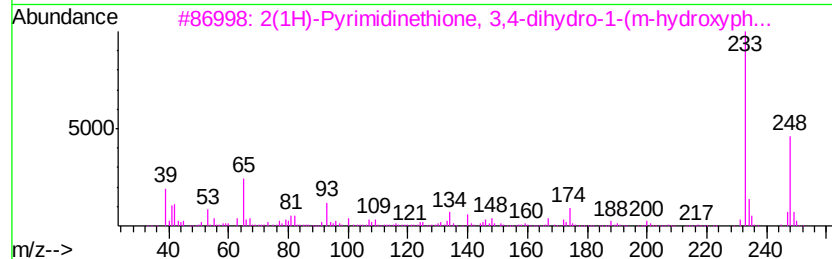
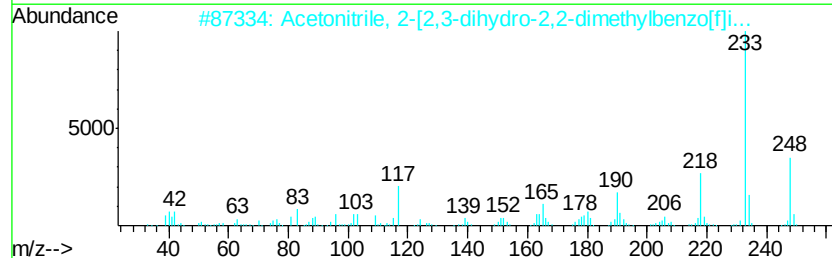
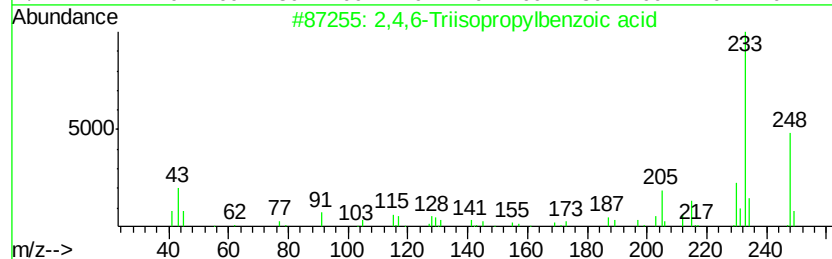
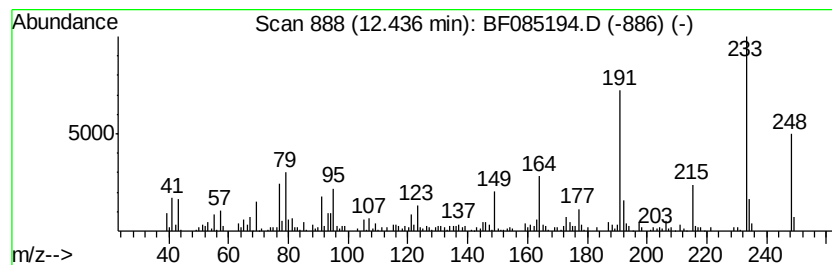
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown12.44 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.44	5.36 ng	314317	Phenanthrene-d10	11.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Triisopropylbenzoic acid	248	C16H24O2	049623-71-4	46
2	Acetonitrile, 2-[2,3-dihydro-2,2...	248	C17H16N2	144498-83-9	46
3	2(1H)-Pyrimidinethione, 3,4-dihy...	248	C13H16N2OS	063704-47-2	43
4	3,5-di-tert-Butyl-4-hydroxyaceto...	248	C16H24O2	014035-33-7	38
5	Phenanthro[9,10-d]oxazole, 2-met...	233	C16H11NO	021639-88-3	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
Data File : BF085194.D
Acq On : 4 Mar 2016 8:50
Operator : SJ/UM
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Instrument :
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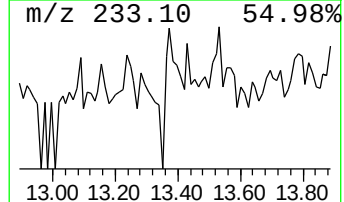
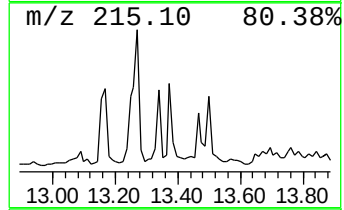
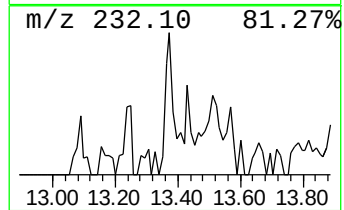
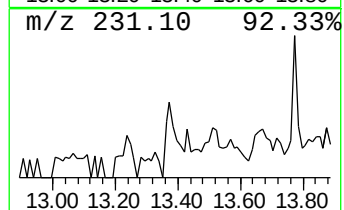
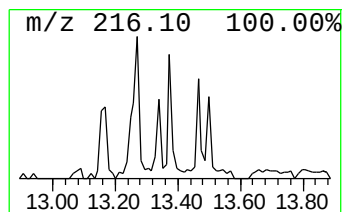
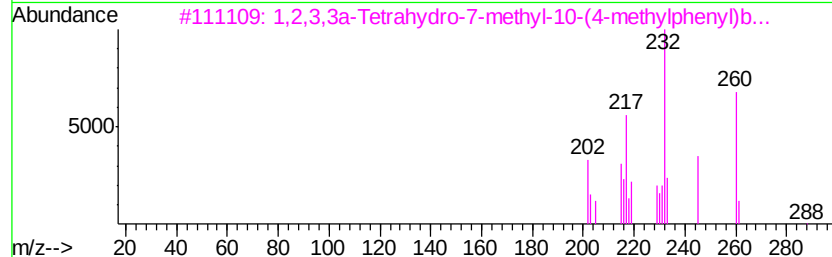
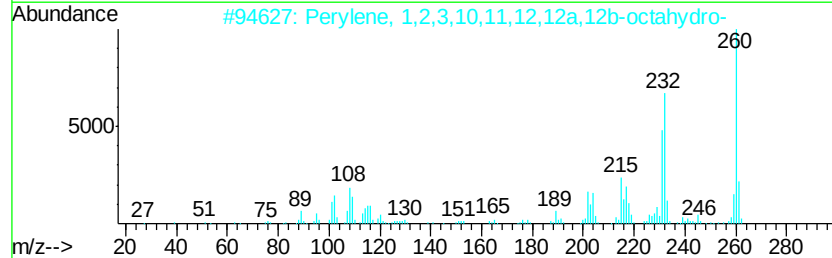
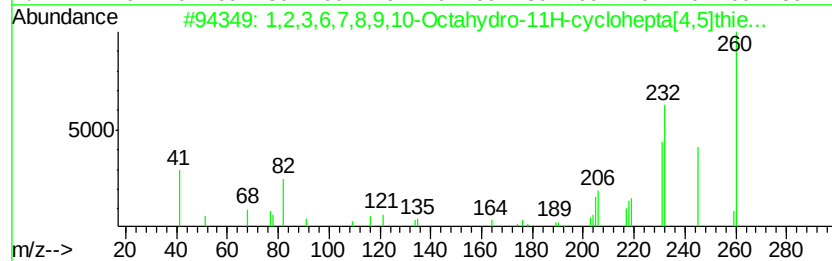
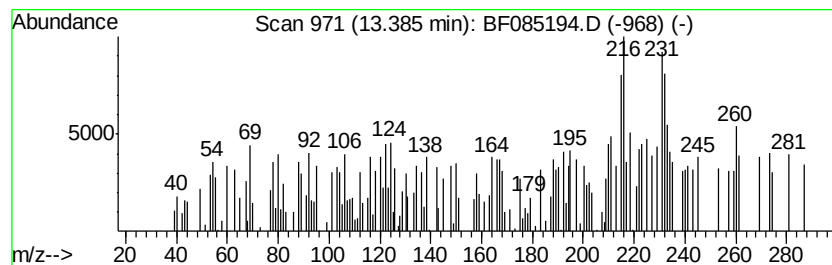
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown13.39 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.39	2.52 ng	126359	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,3,6,7,8,9,10-Octahydro-11H-c...	260	C14H16N2OS	102254-55-7	14
2		Perylene, 1,2,3,10,11,12,12a,12b...	260	C20H20	014930-91-7	14
3		1,2,3,3a-Tetrahydro-7-methyl-10-...	288	C20H20N2	028749-07-7	12
4		Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	11
5		Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
Data File : BF085194.D
Acq On : 4 Mar 2016 8:50
Operator : SJ/UM
Sample : H1649-02 5X
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ALS Vial : 38 Sample Multiplier: 1

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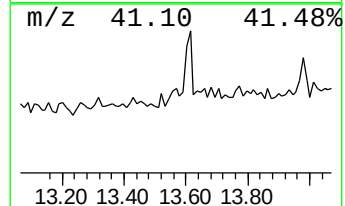
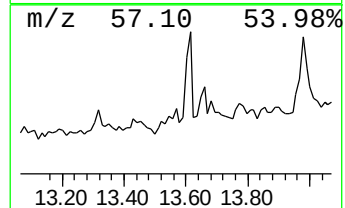
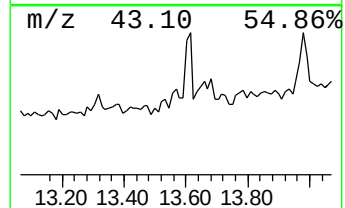
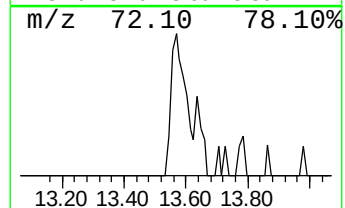
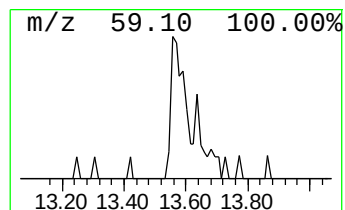
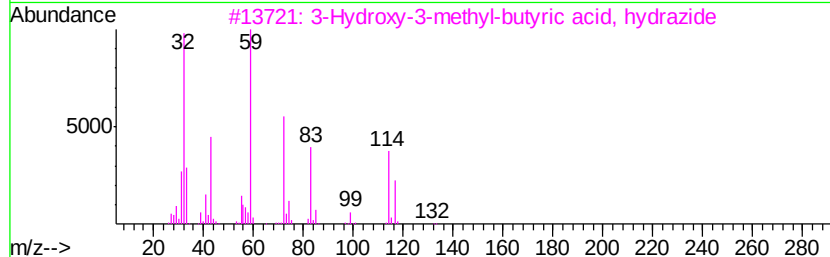
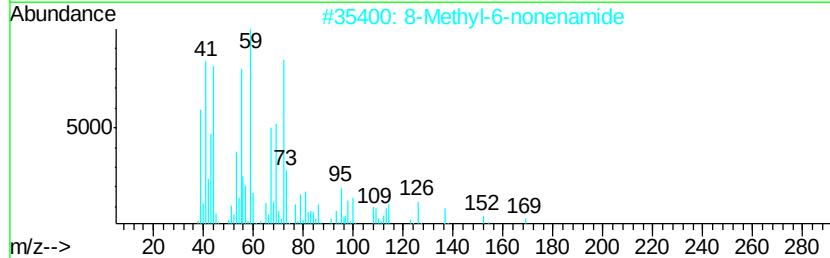
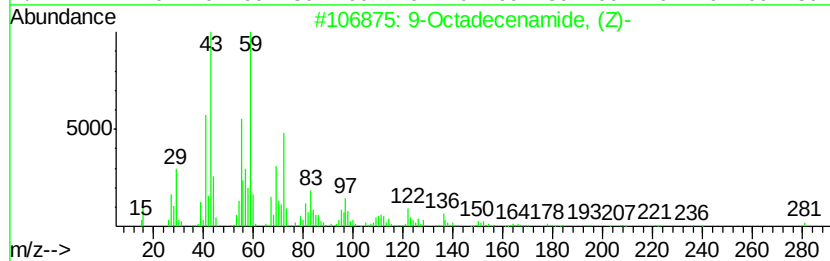
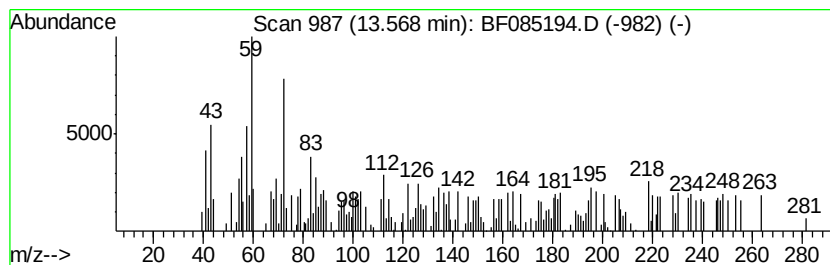
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	2.63 ng	131979	Chrysene-d12	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	60
2		8-Methyl-6-nonenamide	169	C10H19NO	1000293-20-9	49
3		3-Hydroxy-3-methyl-butyric acid,...	132	C5H12N2O2	1000193-80-2	47
4		2-Ethoxyethyl acetate	132	C6H12O3	000111-15-9	35
5		Mannosamine hydrochloride	179	C6H13NO5	1000127-68-8	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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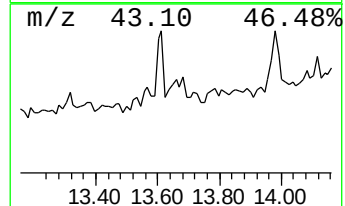
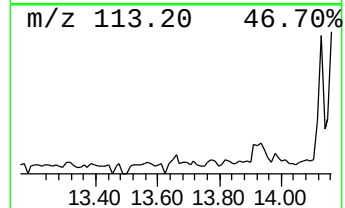
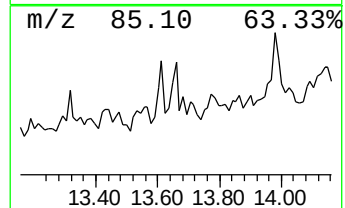
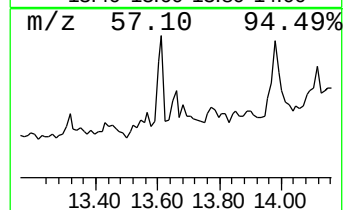
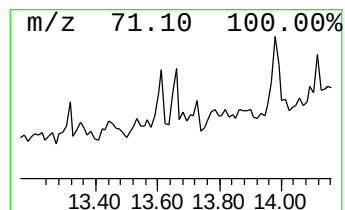
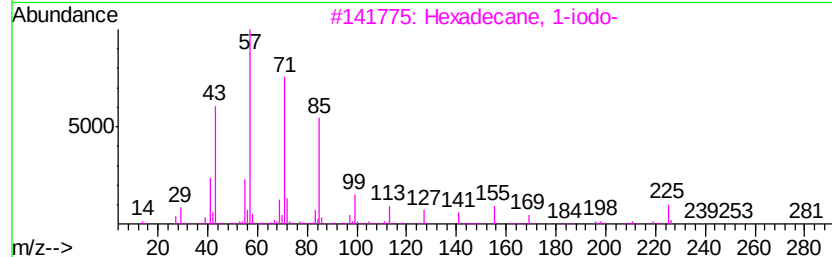
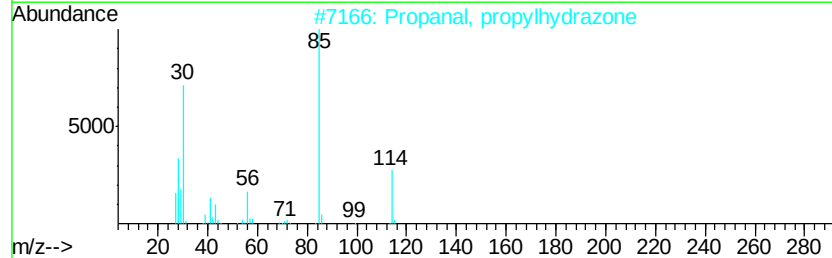
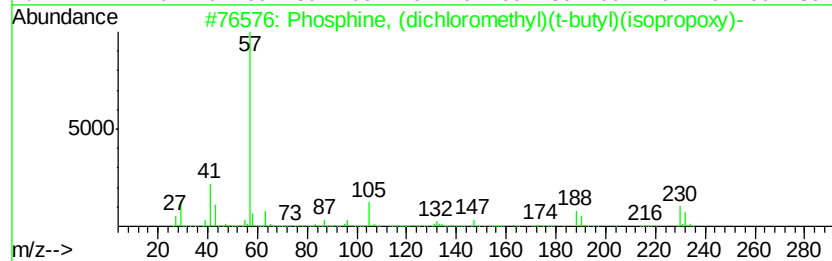
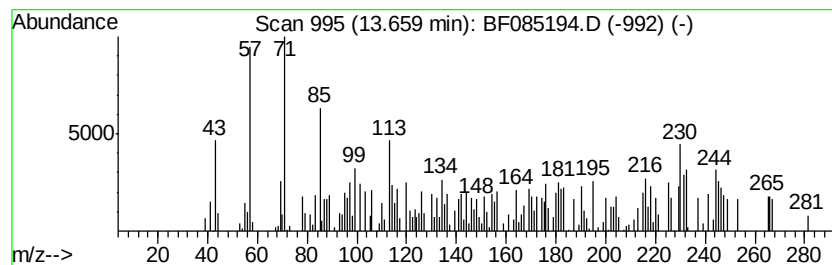
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.66 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.66	2.72 ng	136666	Chrysene-d12	14.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phosphine, (dichloromethyl)(t-bu...	230	C8H17Cl2OP	1000162-70-0	11
2		Propanal, propylhydrazone	114	C6H14N2	019718-39-9	11
3		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	10
4		Pyrrolidine-2-carboxamide, 5-oxo...	246	C8H14N4O3S	1000259-16-7	10
5		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
Data File : BF085194.D
Acq On : 4 Mar 2016 8:50
Operator : SJ/UM
Sample : H1649-02 5X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BT-C3(1)3W-20160301

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.9	ng	219266	1	6.98	894323	20.0
Benzene, 1-ethyl-...	6.50	2.4	ng	105617	1	6.98	894323	20.0
unknown6.74	6.74	22.2	ng	991956	1	6.98	894323	20.0
Benzene, 1,2,3-tr...	6.80	5.4	ng	240180	1	6.98	894323	20.0
Benzene, 1,3-diet...	7.29	2.4	ng	107247	1	6.98	894323	20.0
unknown12.04	12.04	2.7	ng	156150	4	11.50	1172710	20.0
unknown12.44	12.44	5.4	ng	314317	4	11.50	1172710	20.0
unknown13.39	13.39	2.5	ng	126359	5	14.14	1003720	20.0
9-Octadecenamide, ...	13.57	2.6	ng	131979	5	14.14	1003720	20.0
unknown13.66	13.66	2.7	ng	136666	5	14.14	1003720	20.0