

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
 Data File : BF077343.D
 Acq On : 18 Feb 2015 5:31
 Operator : TP/IZ
 Sample : G1302-04
 Misc : W
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FYMW-02A

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-BF021715.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	1.687	65	70	72	rVV2	439048	994746	42.94%	3.194%
2	1.870	83	86	89	rVV	471455	695753	30.03%	2.234%
3	1.915	89	90	93	rVV	38900	46980	2.03%	0.151%
4	2.007	96	98	101	rVB2	40281	63446	2.74%	0.204%
5	2.247	116	119	123	rVB3	52783	136952	5.91%	0.440%
6	2.453	134	137	141	rVB3	63918	142622	6.16%	0.458%
7	2.716	158	160	165	rVB2	61243	113536	4.90%	0.365%
8	2.887	172	175	178	rBV	83859	146849	6.34%	0.472%
9	3.093	191	193	202	rVB3	24278	64221	2.77%	0.206%
10	3.298	204	211	213	rBV2	21145	57799	2.49%	0.186%
11	3.356	213	216	222	rVB	62941	120909	5.22%	0.388%
12	3.470	222	226	232	rBV4	38683	104684	4.52%	0.336%
13	3.630	237	240	244	rVB2	21277	41378	1.79%	0.133%
14	3.710	244	247	249	rBV	16171	35590	1.54%	0.114%
15	3.824	253	257	260	rVV	73284	124339	5.37%	0.399%
16	4.076	276	279	282	rVB2	28209	49767	2.15%	0.160%
17	4.316	297	300	302	rVV3	30734	66405	2.87%	0.213%
18	4.361	302	304	307	rVV	25977	42245	1.82%	0.136%
19	4.430	307	310	312	rVV2	26152	44635	1.93%	0.143%
20	4.487	312	315	318	rVV	31974	59482	2.57%	0.191%
21	4.613	322	326	333	rBV	96019	149702	6.46%	0.481%
22	4.921	350	353	354	rBV2	25795	45698	1.97%	0.147%
23	5.093	362	368	372	rBV6	24456	102497	4.42%	0.329%
24	5.173	373	375	380	rVV	64461	98832	4.27%	0.317%
25	5.253	380	382	386	rVV3	31280	55638	2.40%	0.179%
26	5.504	402	404	407	rBV	624565	702453	30.32%	2.256%
27	5.664	413	418	421	rVB	946861	1719694	74.23%	5.522%
28	5.962	441	444	446	rBV	272796	391577	16.90%	1.257%
29	6.362	476	479	481	rBV	358355	400791	17.30%	1.287%
30	6.716	508	510	512	rVB	751723	673848	29.09%	2.164%
31	6.796	515	517	519	rVV	321952	291887	12.60%	0.937%
32	6.830	519	520	523	rVB	242458	293914	12.69%	0.944%
33	6.887	523	525	526	rBV	424735	420340	18.14%	1.350%
34	6.910	526	527	531	rVB2	240222	342728	14.79%	1.101%

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35	6.990	532	534	537	rVB	485123	437475	18.88%	1.405%
36	7.082	539	542	544	rBV	419544	394303	17.02%	1.266%
37	7.162	547	549	552	rVB	1425526	1285933	55.51%	4.129%
38	7.322	558	563	565	rBV2	60722	119619	5.16%	0.384%
39	7.379	565	568	570	rBV	348170	358209	15.46%	1.150%
40	7.447	570	574	576	rBV	588490	568506	24.54%	1.826%
41	7.562	582	584	585	rBV	833095	765830	33.06%	2.459%
42	7.596	585	587	590	rVV	2456707	2316697	100.00%	7.439%
43	7.687	593	595	597	rBV	276795	250212	10.80%	0.803%
44	7.722	597	598	601	rVB	90450	72258	3.12%	0.232%
45	7.779	601	603	607	rBV	360153	431138	18.61%	1.384%
46	7.870	609	611	614	rBV	152037	197690	8.53%	0.635%
47	7.950	617	618	620	rBV	71981	99956	4.31%	0.321%
48	7.985	620	621	624	rVV2	81944	110549	4.77%	0.355%
49	8.042	624	626	627	rVV	973830	1045108	45.11%	3.356%
50	8.076	627	629	632	rVV2	961276	1362222	58.80%	4.374%
51	8.179	636	638	640	rVV2	36532	49864	2.15%	0.160%
52	8.225	640	642	644	rVB	76472	68209	2.94%	0.219%
53	8.328	649	651	653	rBV	564257	572574	24.72%	1.839%
54	8.362	653	654	656	rVV	158334	122285	5.28%	0.393%
55	8.430	656	660	663	rVB3	41694	83282	3.59%	0.267%
56	8.556	666	671	673	rVV	524712	542854	23.43%	1.743%
57	8.636	675	678	681	rBV	983392	1020628	44.06%	3.277%
58	8.682	681	682	685	rVV2	53619	90013	3.89%	0.289%
59	8.762	685	689	691	rVB3	58181	113073	4.88%	0.363%
60	8.808	691	693	696	rBV	52468	66762	2.88%	0.214%
61	8.956	698	706	707	rVV	605672	721281	31.13%	2.316%
62	8.979	707	708	710	rVV	463958	466781	20.15%	1.499%
63	9.025	710	712	715	rVB	88313	120851	5.22%	0.388%
64	9.276	731	734	736	rVV	102930	184388	7.96%	0.592%
65	9.310	736	737	741	rVV2	55806	74345	3.21%	0.239%
66	9.436	747	748	750	rVB	47595	53566	2.31%	0.172%
67	9.539	755	757	760	rVB	72777	81212	3.51%	0.261%
68	9.653	766	767	769	rVV	53669	68935	2.98%	0.221%
69	9.688	769	770	773	rVB	64503	68705	2.97%	0.221%
70	9.779	777	778	780	rVB2	32727	41730	1.80%	0.134%
71	9.836	780	783	786	rBV2	88114	140188	6.05%	0.450%

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72	9.882	786	787	790	rVV2	30083	47544	2.05%	0.153%
73	9.962	792	794	796	rVV	390324	388691	16.78%	1.248%
74	9.996	796	797	801	rVV3	30534	65447	2.83%	0.210%
75	10.065	801	803	805	rVB2	26470	35248	1.52%	0.113%
76	10.145	807	810	811	rVV3	46539	66142	2.86%	0.212%
77	10.179	811	813	814	rVV2	28183	44661	1.93%	0.143%
78	10.202	814	815	817	rVV	120435	99010	4.27%	0.318%
79	10.305	821	824	826	rVB	1112089	1272645	54.93%	4.087%
80	10.499	839	841	843	rBV	51198	44555	1.92%	0.143%
81	10.602	848	850	853	rBV2	74490	170058	7.34%	0.546%
82	10.659	853	855	858	rVV2	56555	86038	3.71%	0.276%
83	10.705	858	859	861	rVV	29681	46615	2.01%	0.150%
84	10.739	861	862	866	rVB3	45711	60656	2.62%	0.195%
85	10.808	866	868	873	rBV2	98437	166333	7.18%	0.534%
86	10.899	873	876	879	rBV3	30188	69972	3.02%	0.225%
87	11.128	894	896	899	rVV	502191	589665	25.45%	1.894%
88	12.031	973	975	980	rVB	164130	155107	6.70%	0.498%
89	12.111	980	982	984	rBV	260458	276600	11.94%	0.888%
90	12.979	1055	1058	1060	rBV	613927	615968	26.59%	1.978%
91	13.688	1117	1120	1122	rBV	105304	124307	5.37%	0.399%
92	14.328	1174	1176	1180	rBV	44360	60438	2.61%	0.194%
93	14.671	1204	1206	1209	rBV	33202	46151	1.99%	0.148%
94	14.980	1230	1233	1235	rBV	1099510	1328908	57.36%	4.267%
95	15.563	1282	1284	1287	rBV	37008	42162	1.82%	0.135%
96	16.145	1332	1335	1344	rBV	314494	496548	21.43%	1.595%
97	16.271	1344	1346	1349	rBV	571783	620633	26.79%	1.993%
98	16.968	1405	1407	1414	rBV3	21397	43437	1.87%	0.139%
99	18.031	1497	1500	1503	rBV	474636	593611	25.62%	1.906%
100	18.671	1554	1556	1560	rBV	35179	74958	3.24%	0.241%

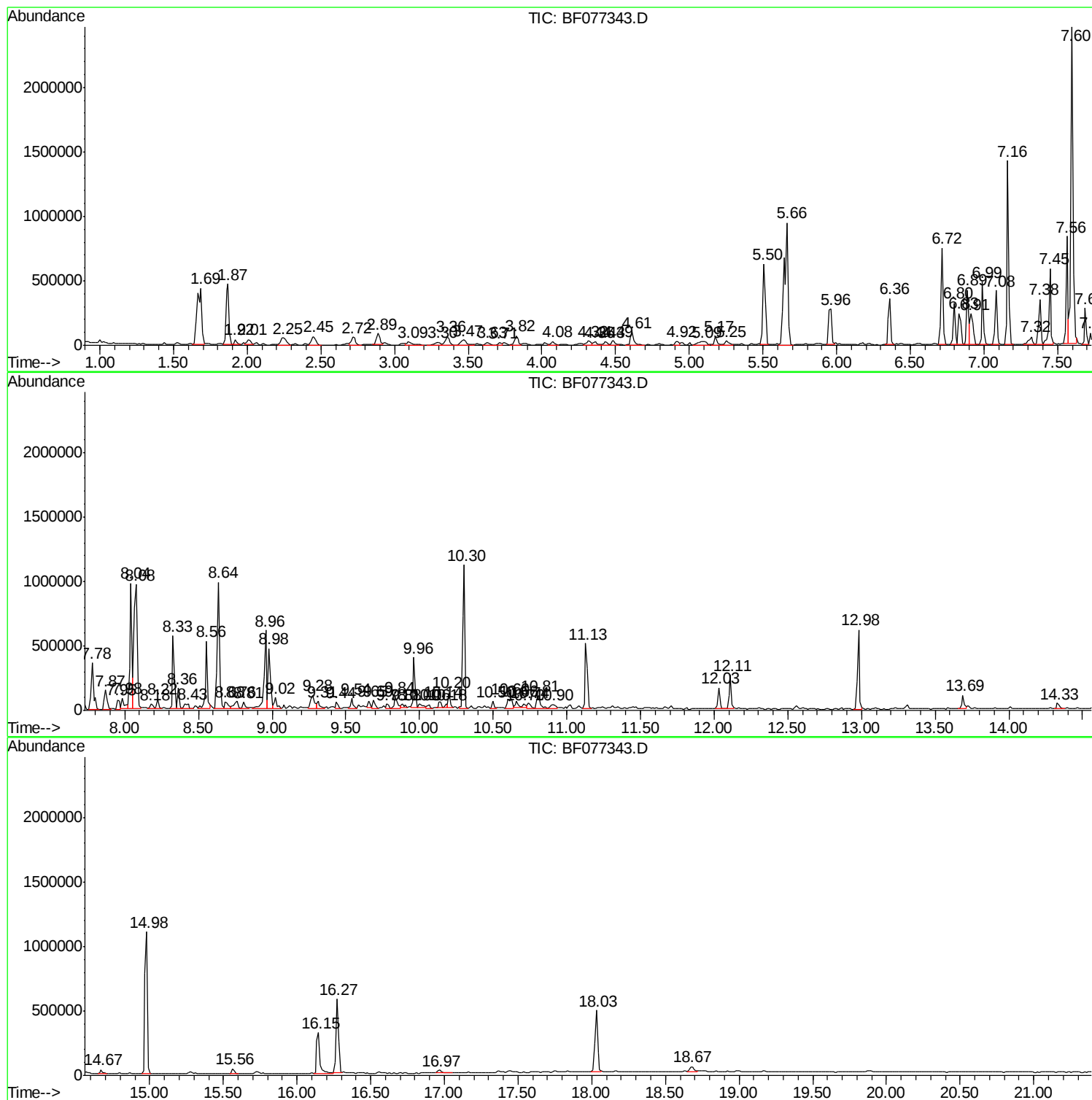
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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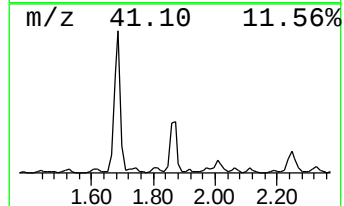
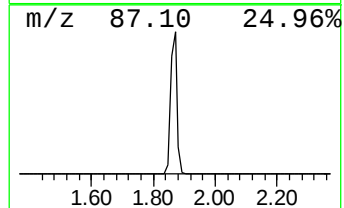
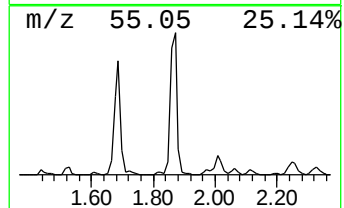
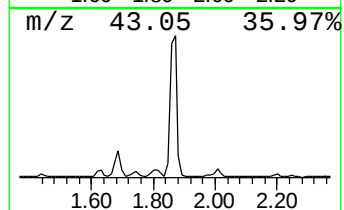
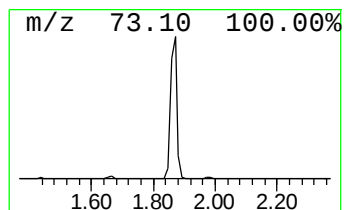
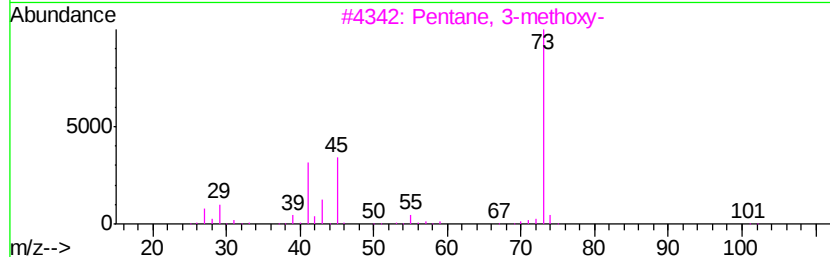
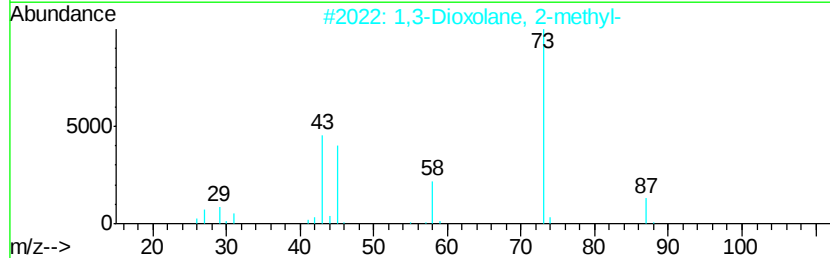
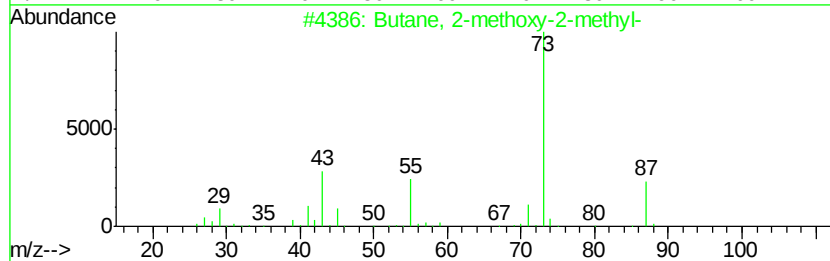
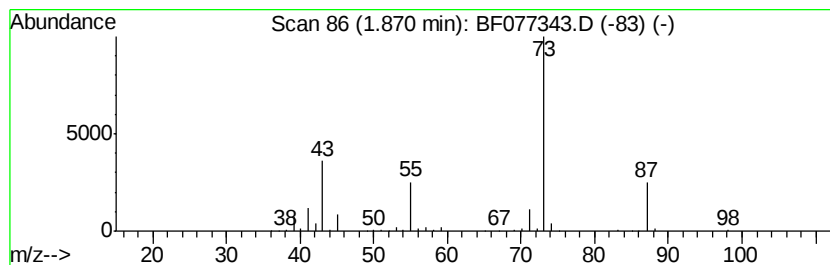
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.87	38.85 ng	695753	1,4-Dichlorobenzene-d4	7.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10
5			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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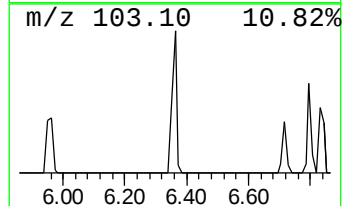
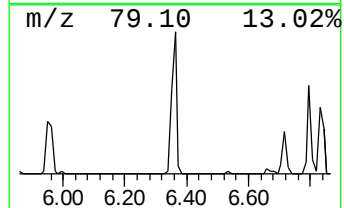
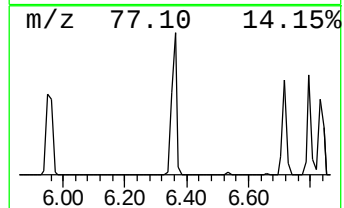
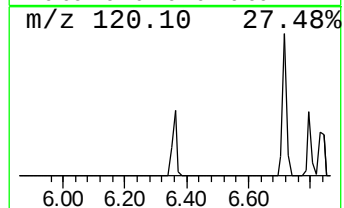
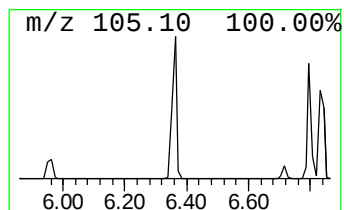
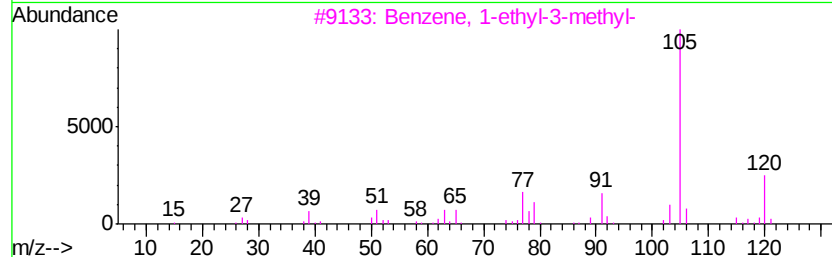
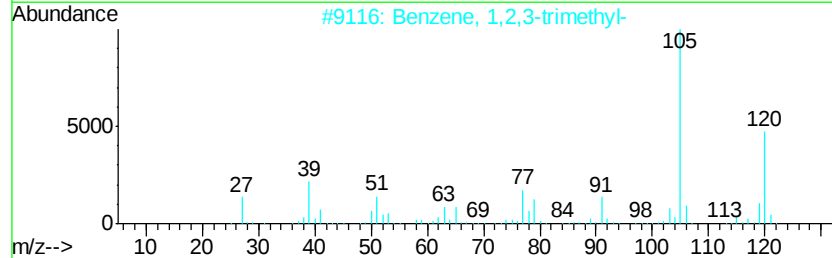
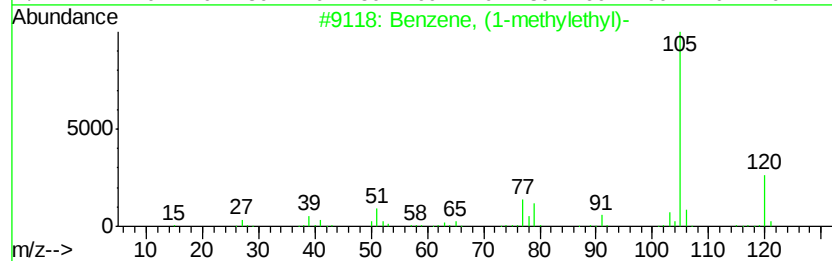
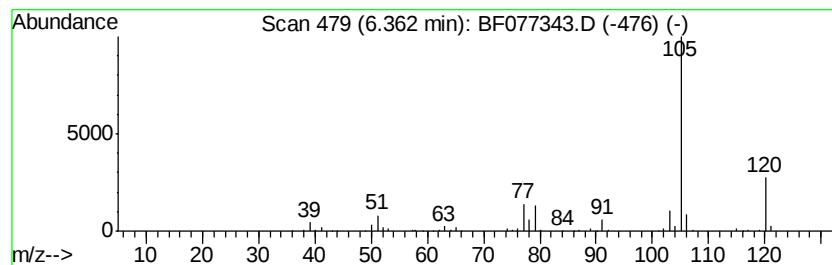
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, (1-methylethyl)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.36	22.38 ng	400791	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	86
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	80
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	78



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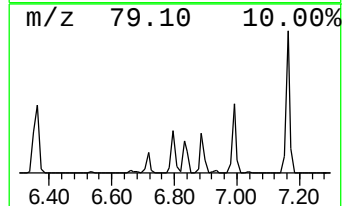
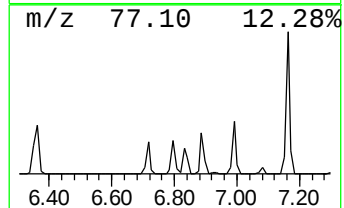
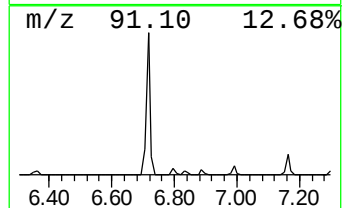
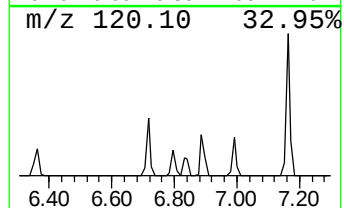
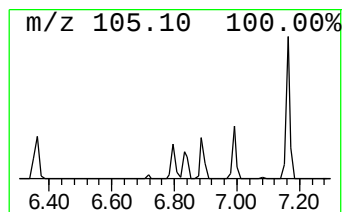
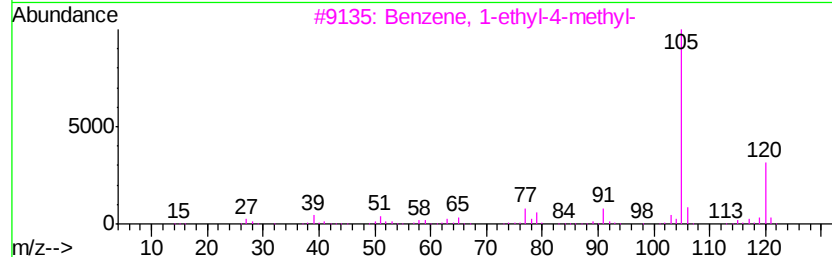
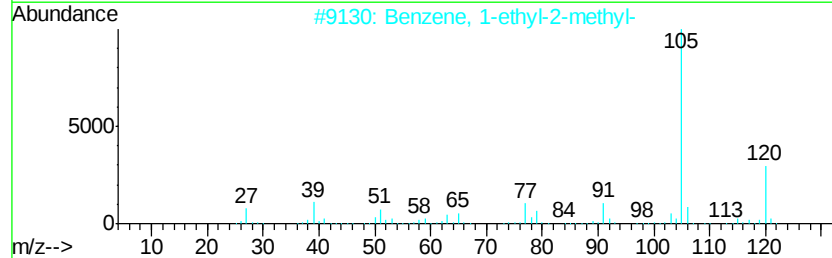
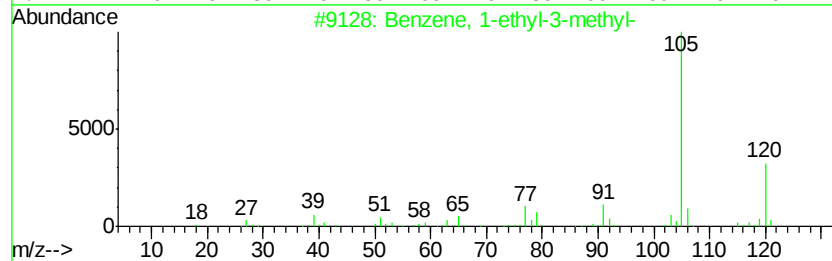
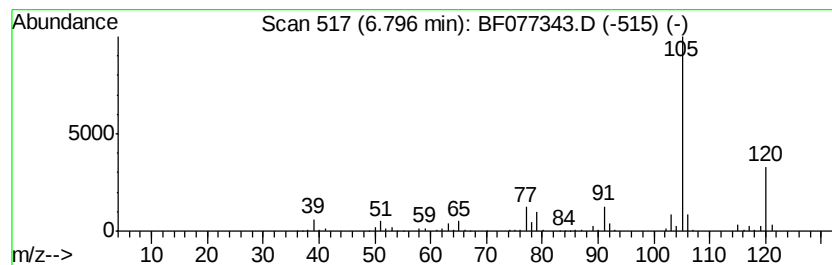
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	16.30 ng	291887	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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BNA_F
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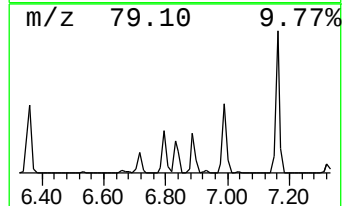
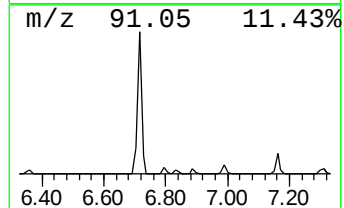
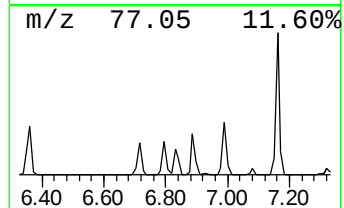
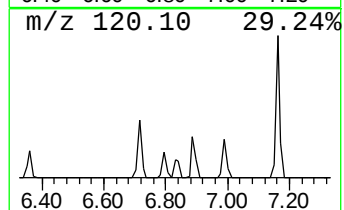
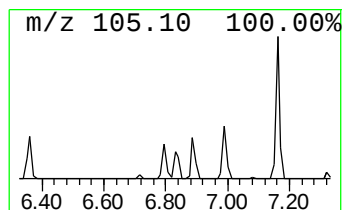
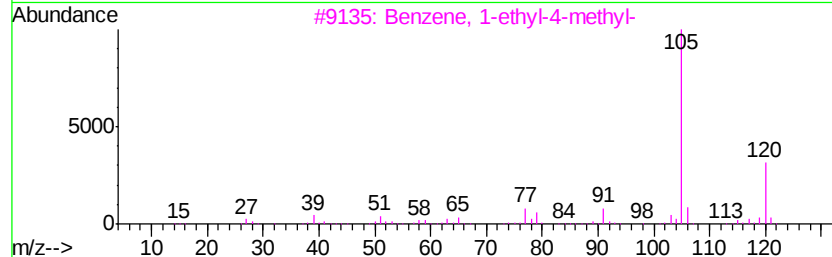
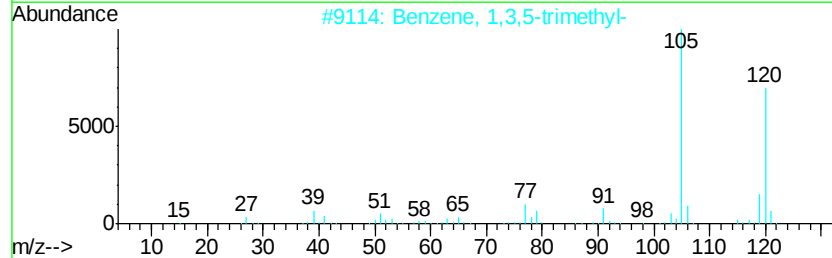
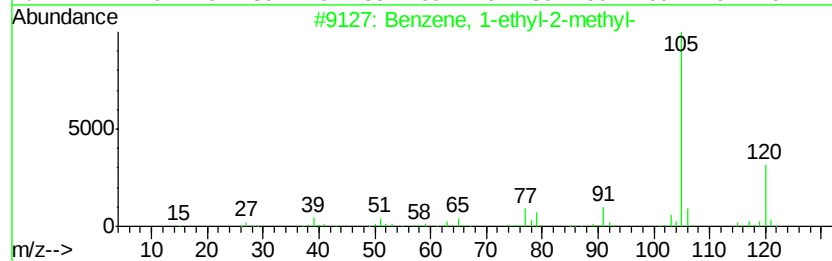
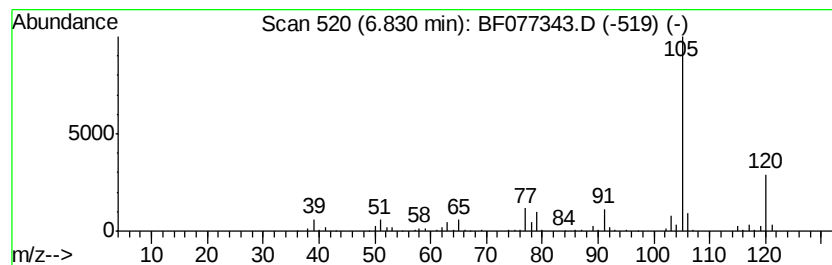
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.83	16.41 ng	293914	1,4-Dichlorobenzene-d4	7.38

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
Data File : BF077343.D
Acq On : 18 Feb 2015 5:31
Operator : TP/IZ
Sample : G1302-04
Misc : W
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FYMW-02A

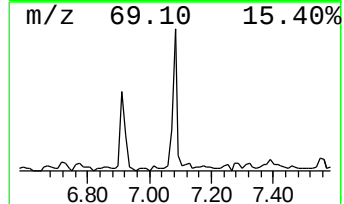
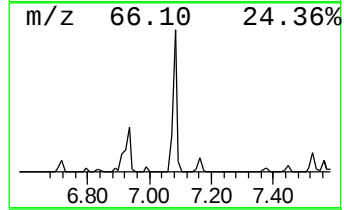
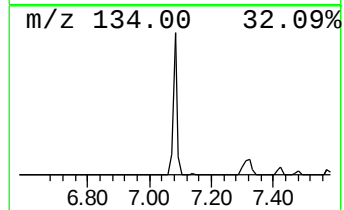
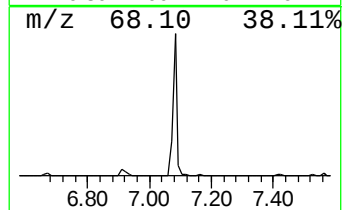
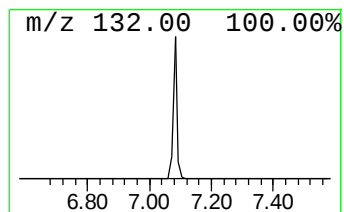
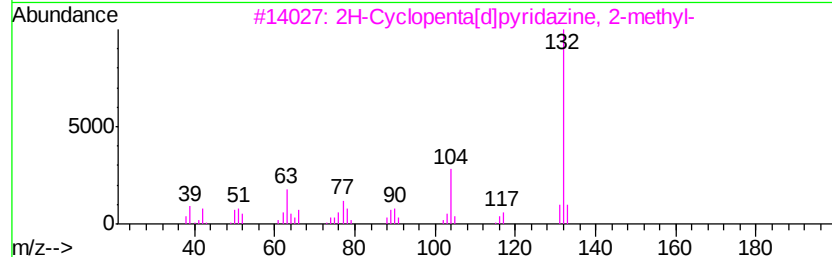
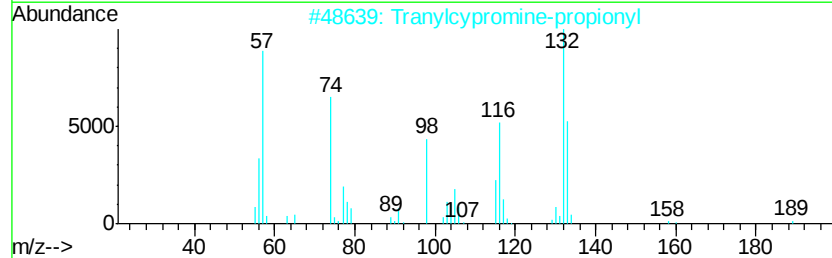
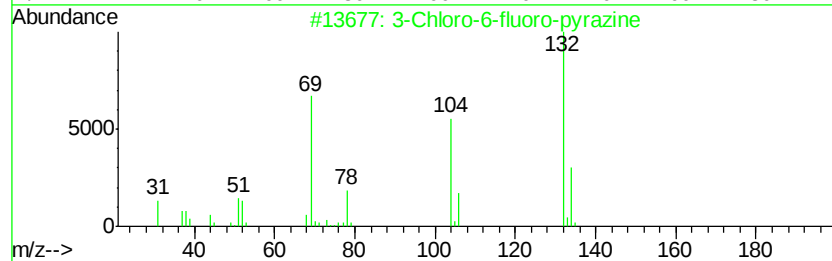
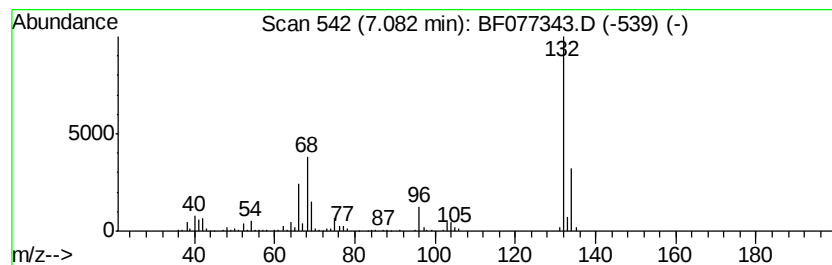
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown7.08 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	22.02 ng	394303	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		5-Aminoindole	132	C8H8N2	005192-03-0	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
Data File : BF077343.D
Acq On : 18 Feb 2015 5:31
Operator : TP/IZ
Sample : G1302-04
Misc : W
ALS Vial : 23 Sample Multiplier: 1

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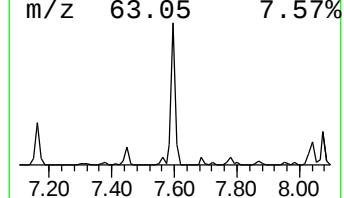
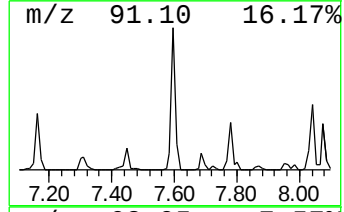
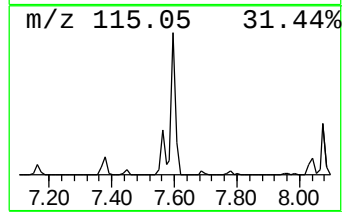
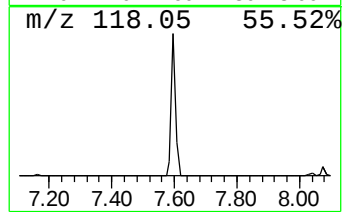
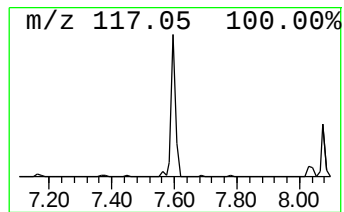
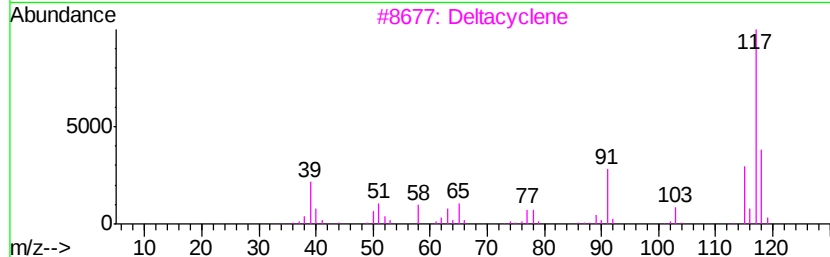
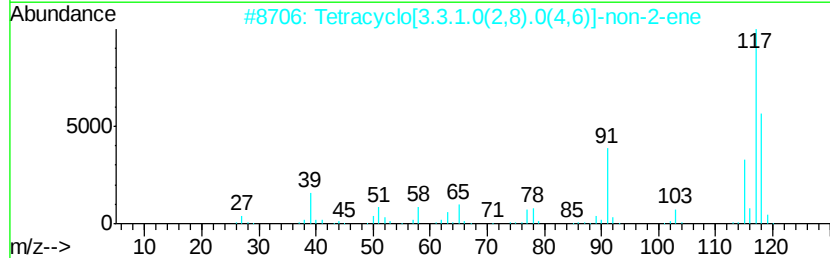
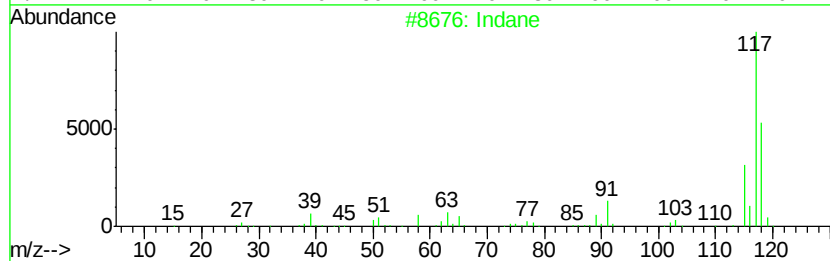
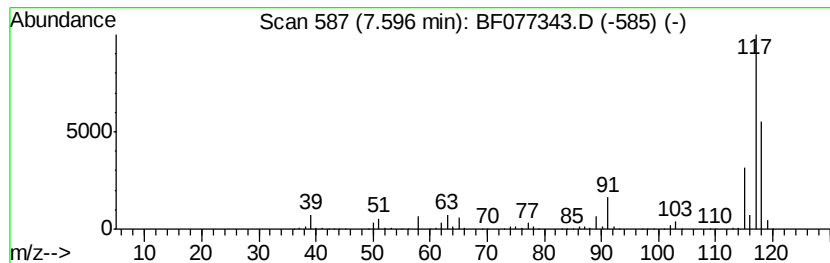
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	129.35 ng	2316700	1,4-Dichlorobenzene-d4	7.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	93
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	83
3			Deltacyclene	118	C9H10	007785-10-6	72
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	72
5			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
Data File : BF077343.D
Acq On : 18 Feb 2015 5:31
Operator : TP/IZ
Sample : G1302-04
Misc : W
ALS Vial : 23 Sample Multiplier: 1

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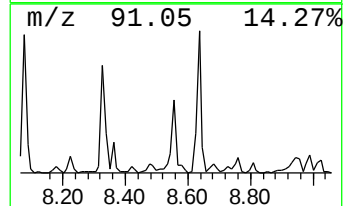
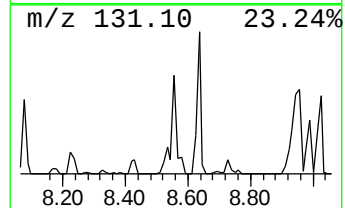
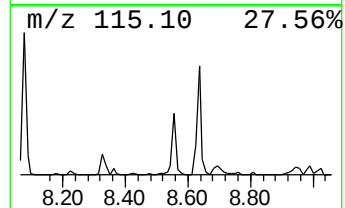
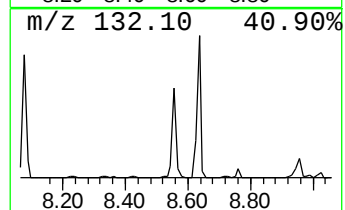
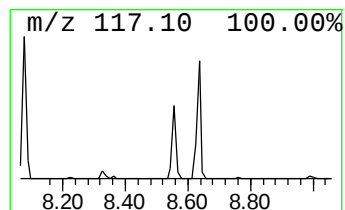
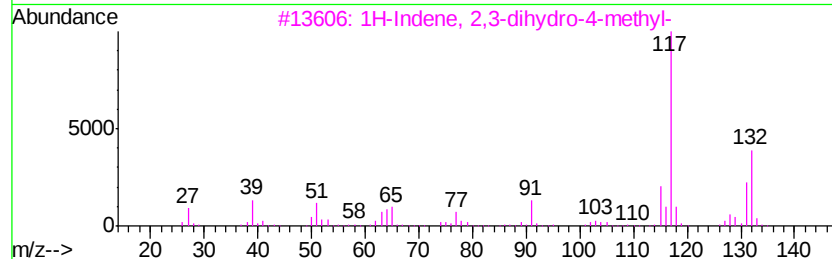
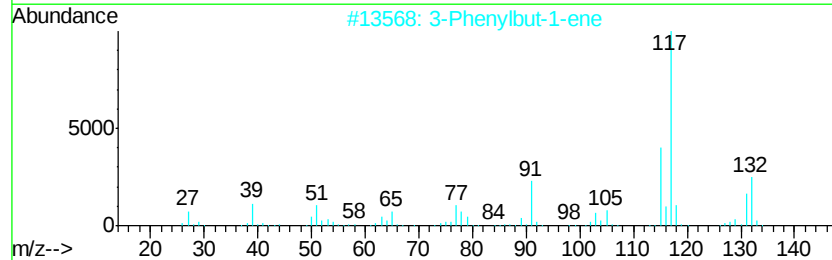
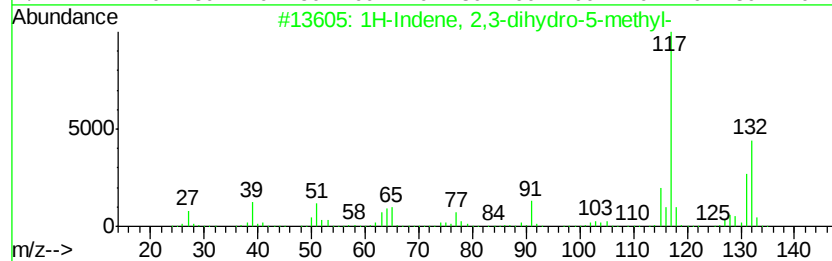
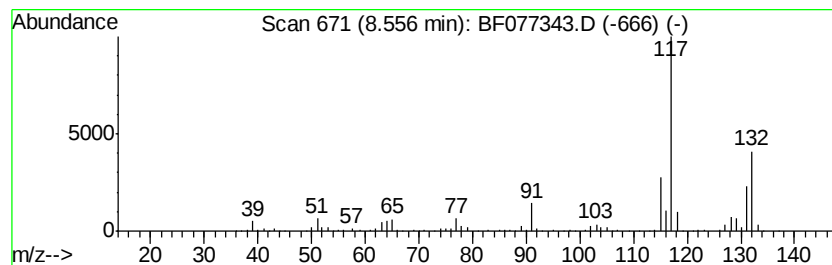
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	15.05 ng	542854	Napthalene-d8	8.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
Data File : BF077343.D
Acq On : 18 Feb 2015 5:31
Operator : TP/IZ
Sample : G1302-04
Misc : W
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FYMW-02A

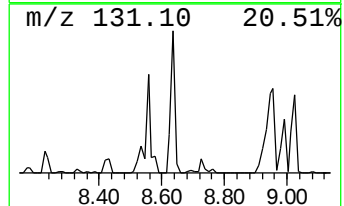
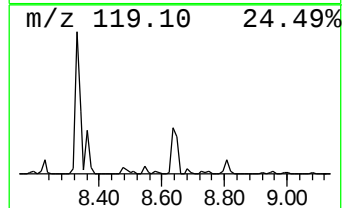
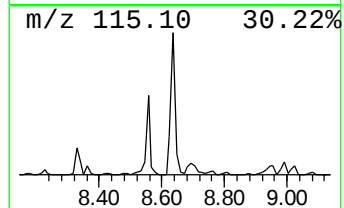
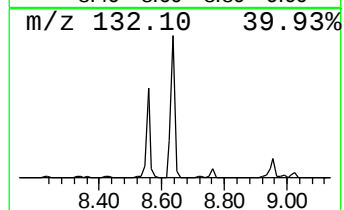
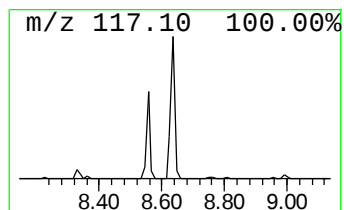
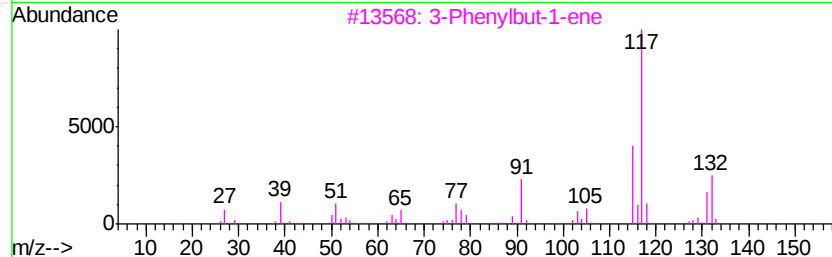
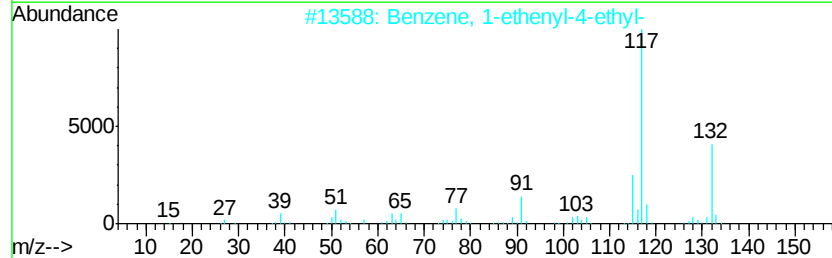
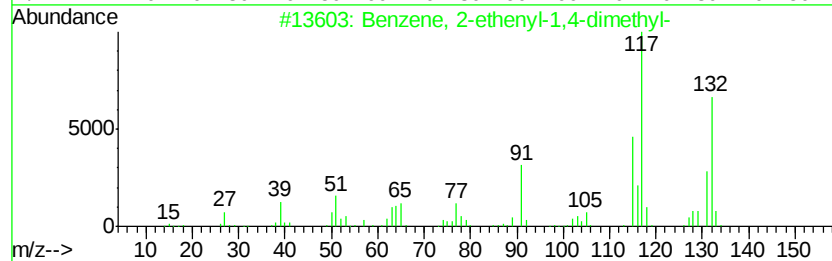
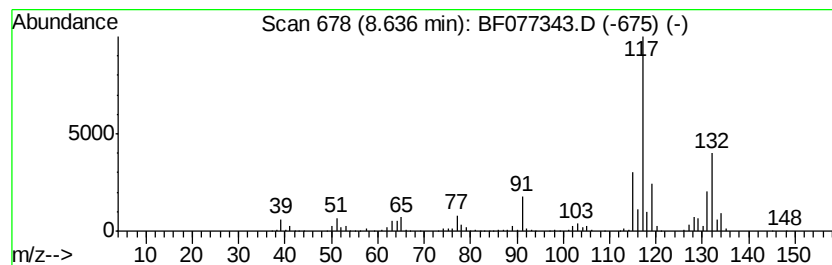
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.64	28.30 ng	1020630	Naphthalene-d8	8.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	83
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
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Operator : TP/IZ
Sample : G1302-04
Misc : W
ALS Vial : 23 Sample Multiplier: 1

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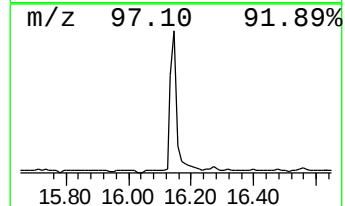
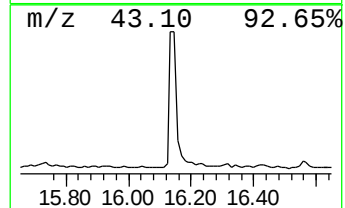
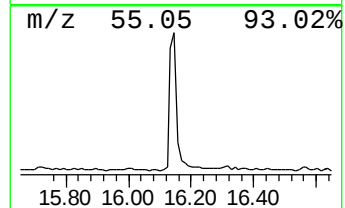
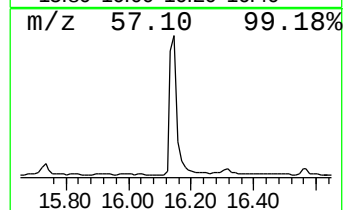
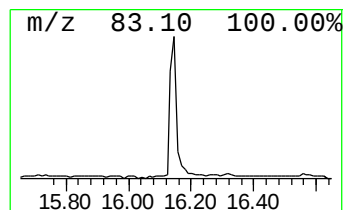
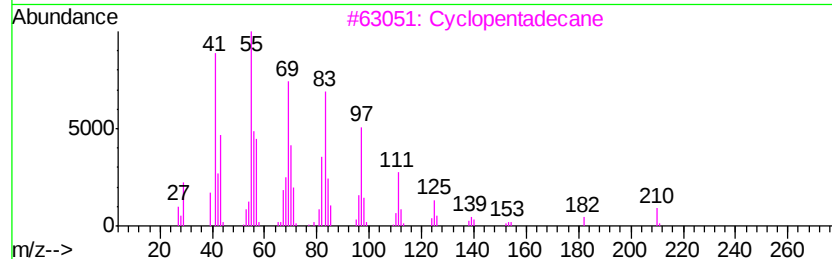
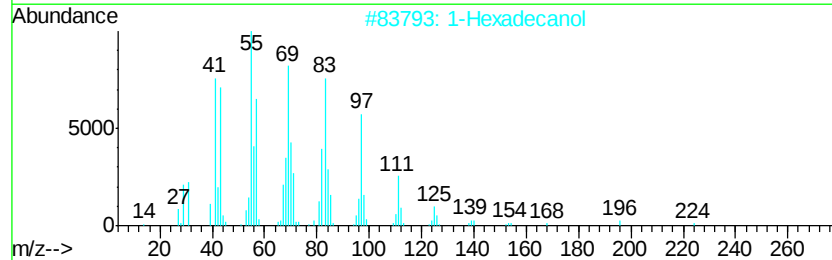
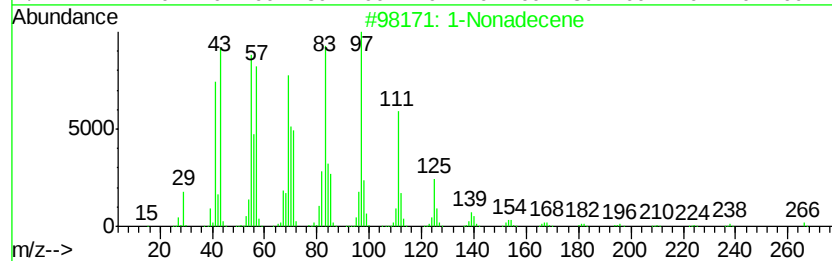
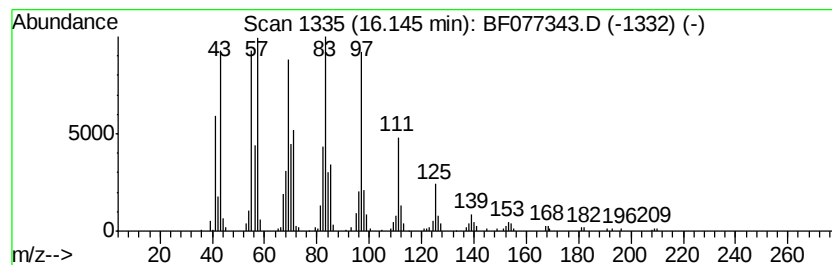
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 1-Nonadecene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	16.00 ng	496548	Chrysene-d12	16.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	91
2		1-Hexadecanol	242	C16H34O	036653-82-4	91
3		Cyclopentadecane	210	C15H30	000295-48-7	91
4		Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	90
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021715\
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Acq On : 18 Feb 2015 5:31
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Sample : G1302-04
Misc : W
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021715.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
Butane, 2-methoxy...	1.87	38.9	ng	695753	1	7.38	358209	20.0
Benzene, (1-methy...	6.36	22.4	ng	400791	1	7.38	358209	20.0
Benzene, 1-ethyl-...	6.80	16.3	ng	291887	1	7.38	358209	20.0
Benzene, 1-ethyl-...	6.83	16.4	ng	293914	1	7.38	358209	20.0
unknown7.08	7.08	22.0	ng	394303	1	7.38	358209	20.0
Indane	7.60	129.3	ng	2316700	1	7.38	358209	20.0
1H-Indene, 2,3-di...	8.56	15.1	ng	542854	2	8.96	721281	20.0
Benzene, 2-etheny...	8.64	28.3	ng	1020630	2	8.96	721281	20.0
1-Nonadecene	16.15	16.0	ng	496548	5	16.27	620633	20.0