

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
 Data File : BF087693.D
 Acq On : 5 Jun 2016 7:11
 Operator : UM/SJ
 Sample : H3344-05
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-18

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-BF060416.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	1.770	15	16	20	rVV	274484	338047	7.97%	0.783%
2	1.838	20	22	26	rVB	210571	263601	6.22%	0.610%
3	1.907	26	28	47	rVB	1927580	3682662	86.86%	8.527%
4	2.147	47	49	68	rVB	78504	250824	5.92%	0.581%
5	3.827	193	196	199	rBV	156221	218916	5.16%	0.507%
6	5.039	299	302	306	rBV	146953	181906	4.29%	0.421%
7	5.313	324	326	329	rBV	50610	58476	1.38%	0.135%
8	5.450	335	338	340	rVB	1910202	1864979	43.99%	4.318%
9	5.690	357	359	362	rVB	790236	694681	16.39%	1.609%
10	6.022	386	388	390	rVB	113355	91733	2.16%	0.212%
11	6.479	424	428	430	rBV	1344811	1226503	28.93%	2.840%
12	6.547	432	434	435	rBV	131686	146110	3.45%	0.338%
13	6.627	438	441	443	rBV	2808508	3245003	76.54%	7.514%
14	6.685	445	446	449	rVB2	141291	138256	3.26%	0.320%
15	6.856	459	461	465	rBV	857407	761740	17.97%	1.764%
16	6.913	465	466	471	rVB2	266478	309246	7.29%	0.716%
17	7.016	472	475	476	rVV	2852336	2746763	64.79%	6.360%
18	7.039	476	477	479	rVB	426508	351320	8.29%	0.813%
19	7.119	482	484	487	rBV	2274381	2341641	55.23%	5.422%
20	7.427	507	511	513	rVB	2177639	2763958	65.19%	6.400%
21	7.553	520	522	524	rBV2	25462	45052	1.06%	0.104%
22	7.656	530	531	534	rVB2	49971	54776	1.29%	0.127%
23	7.896	548	552	553	rBV3	330768	488958	11.53%	1.132%
24	7.930	553	555	558	rVB	414710	696379	16.43%	1.612%
25	8.170	571	576	578	rBV2	2037014	3182213	75.06%	7.368%
26	8.216	578	580	582	rVV	260595	220005	5.19%	0.509%
27	8.296	585	587	592	rBV	69181	77908	1.84%	0.180%
28	8.490	600	604	607	rBV	24687	47714	1.13%	0.110%
29	8.616	613	615	616	rVV	50119	72771	1.72%	0.169%
30	8.639	616	617	619	rVV	190649	195196	4.60%	0.452%
31	8.730	623	625	627	rVB	50318	43741	1.03%	0.101%
32	8.810	630	632	634	rVB	81229	76931	1.81%	0.178%
33	8.959	642	645	647	rBV	722307	740337	17.46%	1.714%
34	9.222	665	668	670	rVV	3163917	4239559	100.00%	9.817%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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35	9.325	673	677	678 rBV	70053	75461	1.78%	0.175%
36	9.428	681	686	688 rVB3	23434	60187	1.42%	0.139%
37	9.553	696	697	699 rVB	52128	53745	1.27%	0.124%
38	9.611	699	702	706 rVV2	24080	45802	1.08%	0.106%
39	9.668	706	707	712 rVB	26795	47849	1.13%	0.111%
40	9.759	712	715	717 rVB	477384	424042	10.00%	0.982%
41	9.896	725	727	729 rBV	1210144	1124343	26.52%	2.603%
42	9.931	729	730	731 rVB	191091	131088	3.09%	0.304%
43	10.136	746	748	750 rBV	75413	69101	1.63%	0.160%
44	10.685	793	796	799 rBV	2246731	2519986	59.44%	5.835%
45	10.822	804	808	812 rVB2	52914	93265	2.20%	0.216%
46	11.291	847	849	854 rVB2	27610	52059	1.23%	0.121%
47	11.382	854	857	859 rBV	1088363	1145138	27.01%	2.652%
48	11.908	901	903	906 rBV	76133	93328	2.20%	0.216%
49	12.697	970	972	976 rBV	69977	84994	2.00%	0.197%
50	12.971	993	996	998 rBV	3330412	3531407	83.30%	8.177%
51	14.011	1084	1087	1089 rBV	1135372	1029432	24.28%	2.384%
52	15.440	1209	1212	1215 rBV	620063	747878	17.64%	1.732%

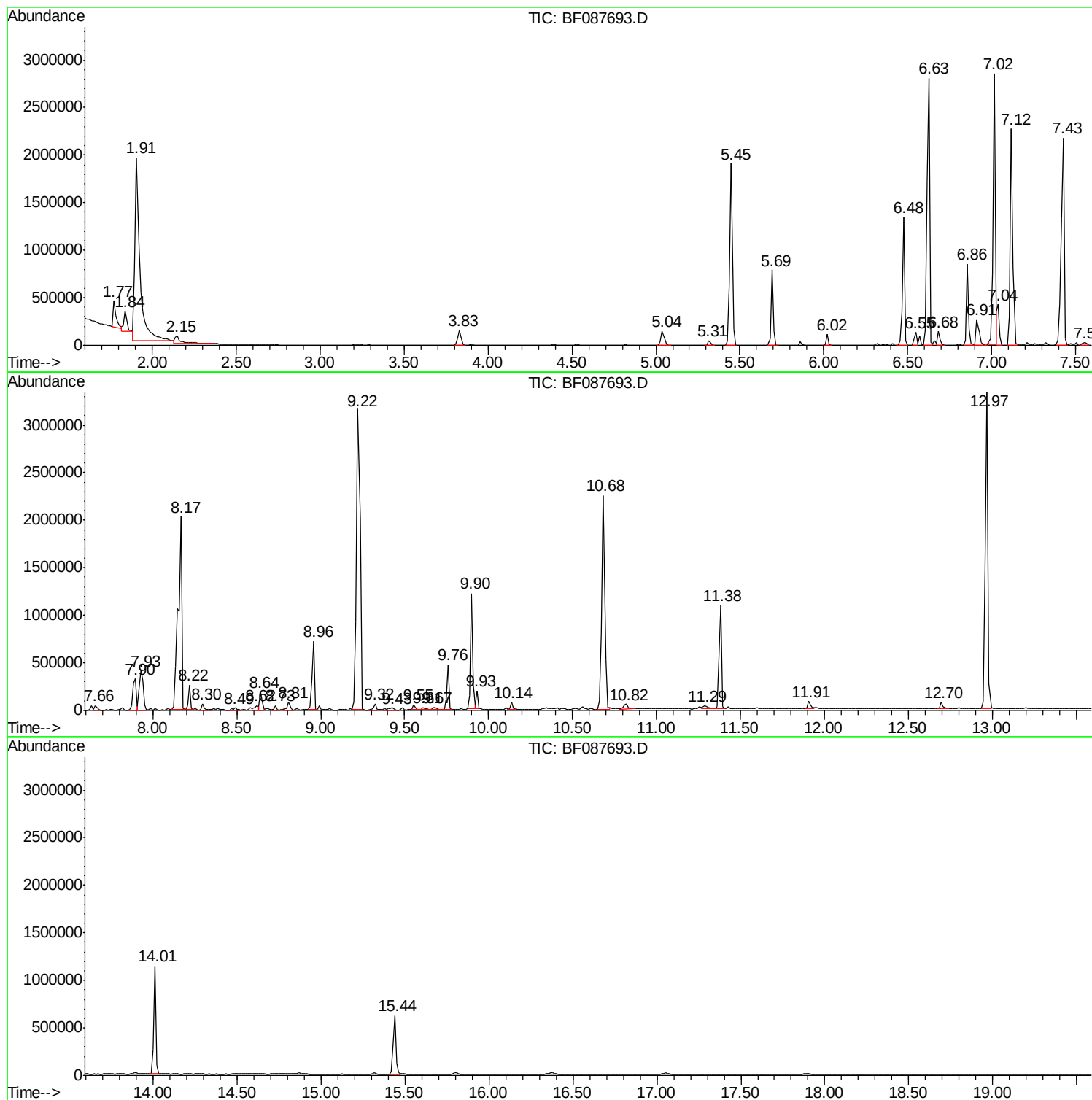
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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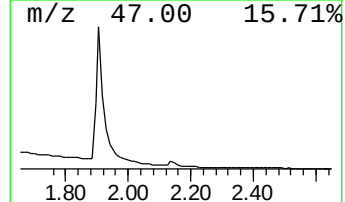
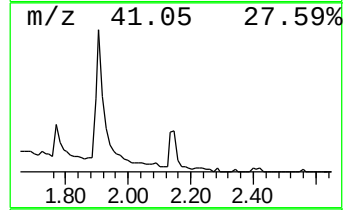
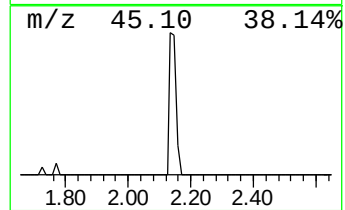
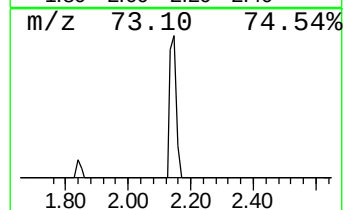
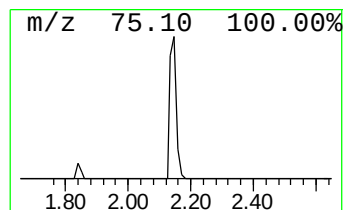
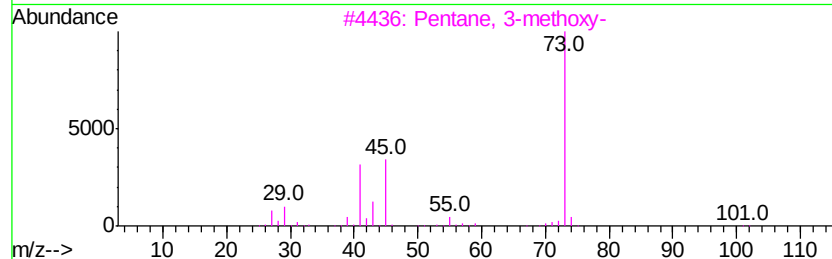
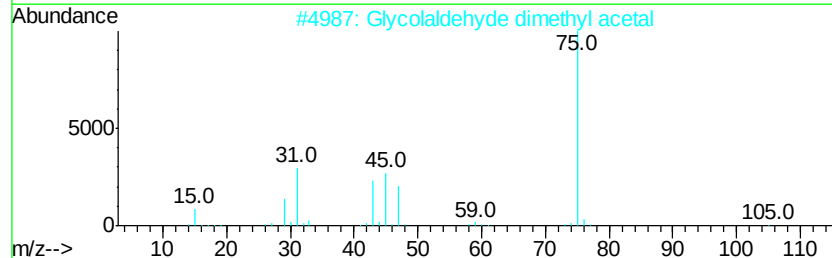
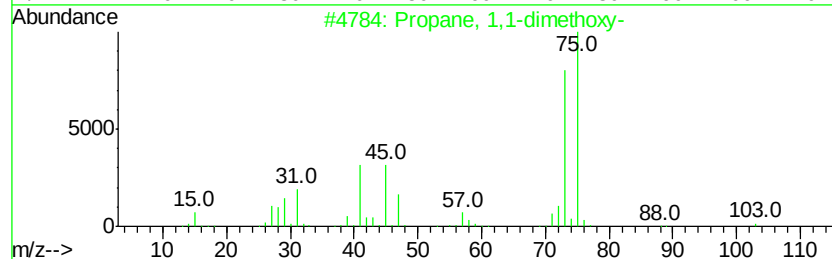
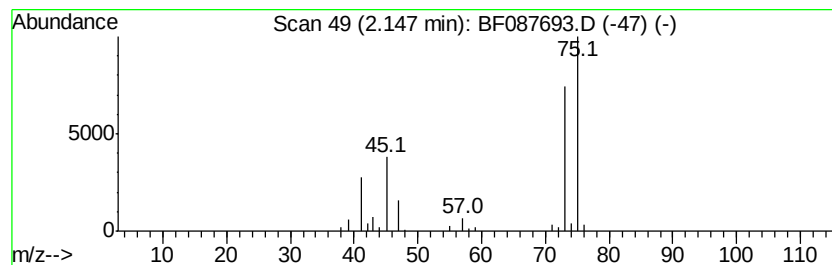
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Propane, 1,1-dimethoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	6.59 ng	250824	1,4-Dichlorobenzene-d4	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2		Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	36
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	16
4		1,3-Dioxolane, 2-(1-bromoethyl)-	180	C5H9BrO2	005267-73-2	16
5		Glycine	75	C2H5NO2	000056-40-6	9



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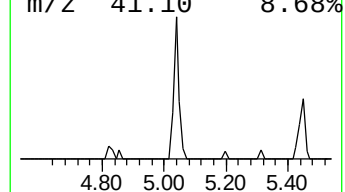
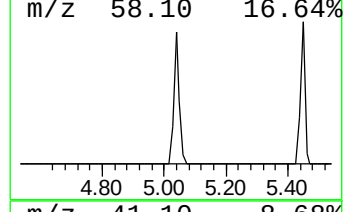
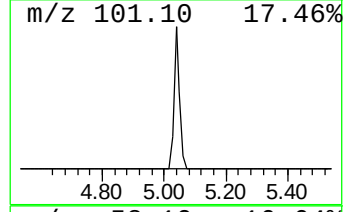
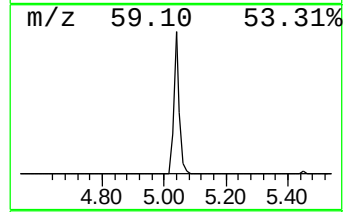
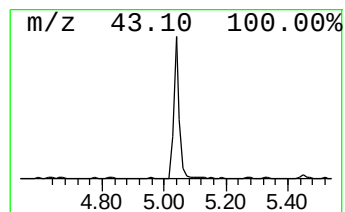
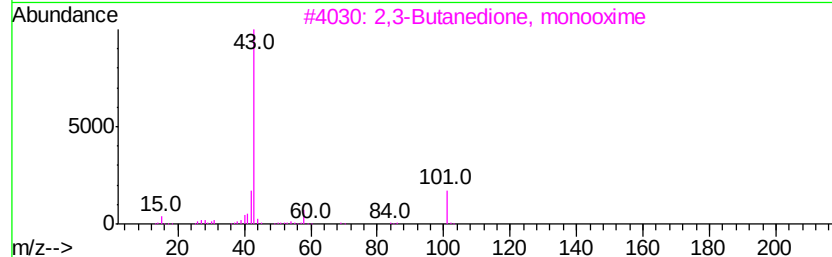
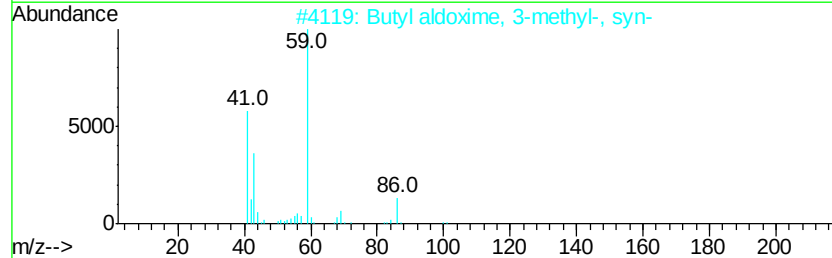
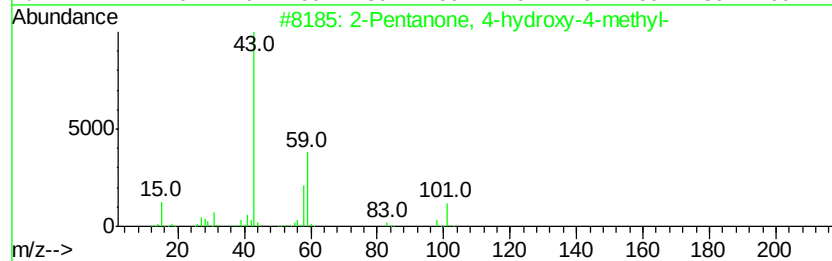
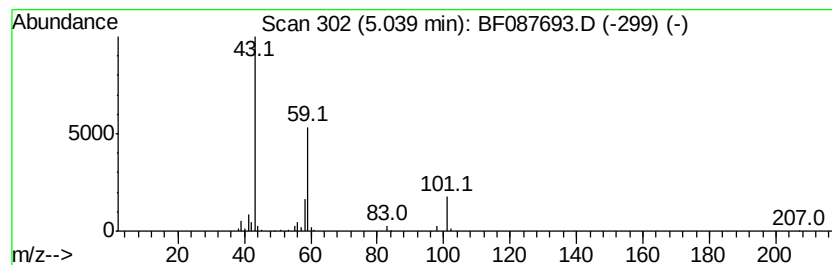
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Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	4.78 ng	181906	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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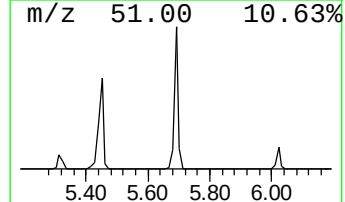
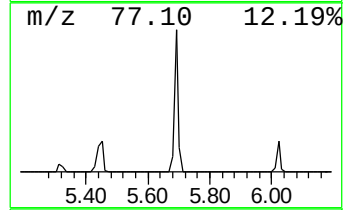
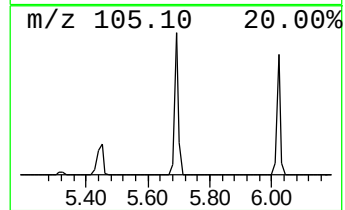
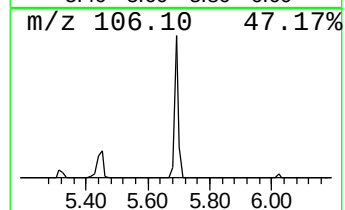
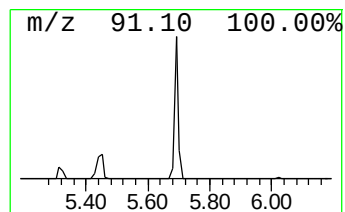
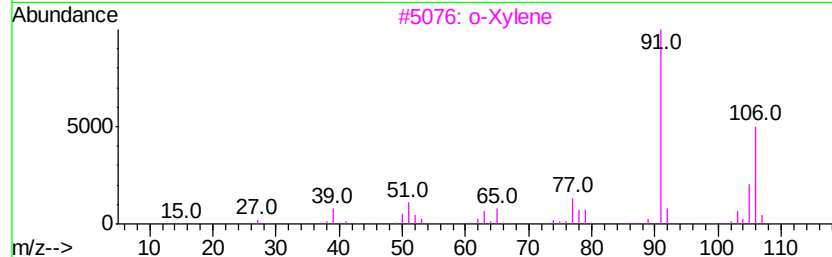
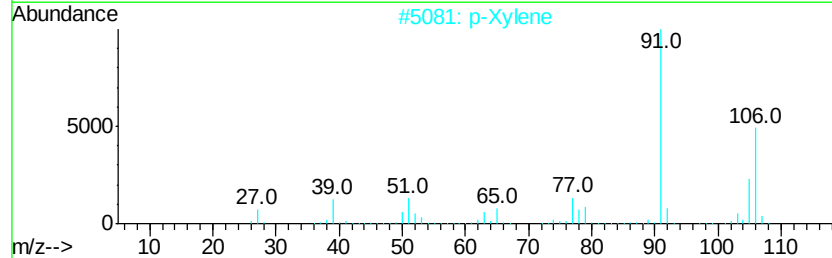
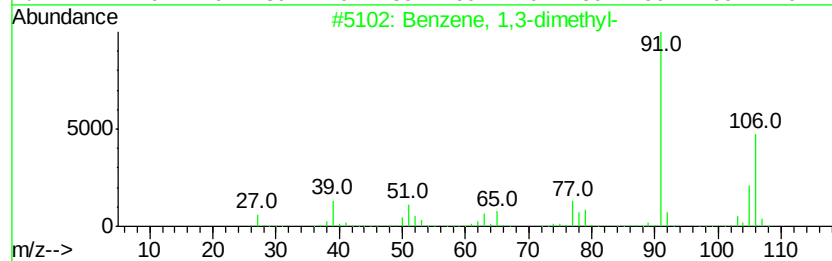
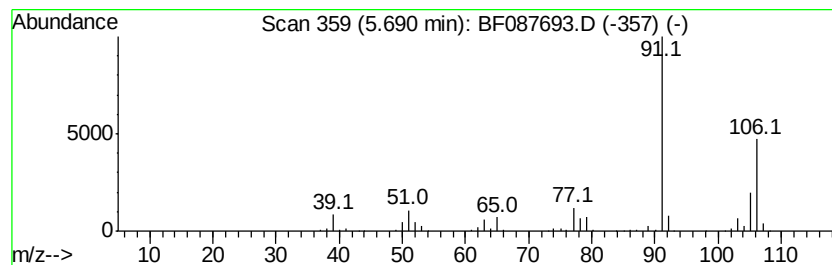
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.69	18.24 ng	694681	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	Ethylbenzene	106	C8H10	000100-41-4	87
5	Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	86



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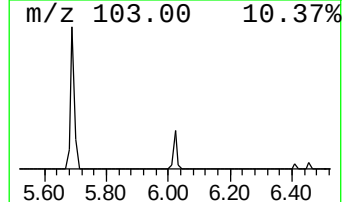
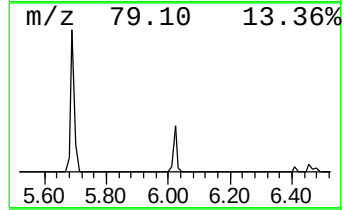
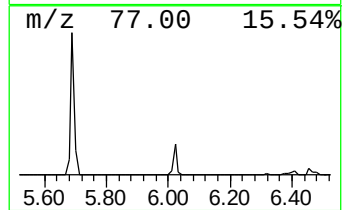
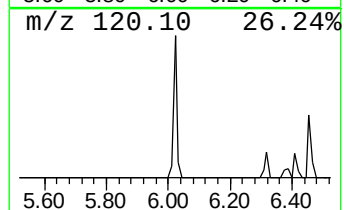
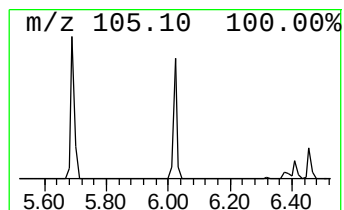
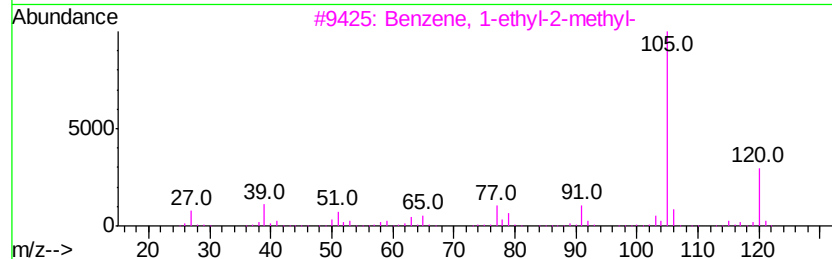
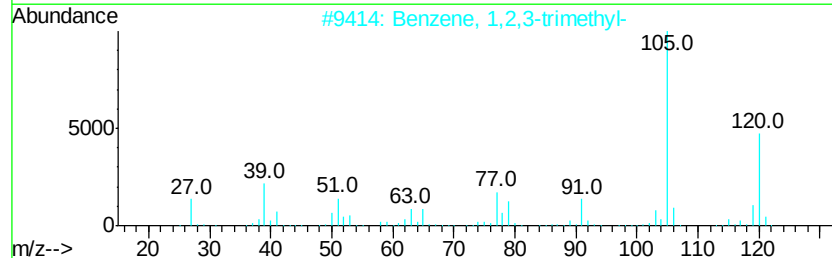
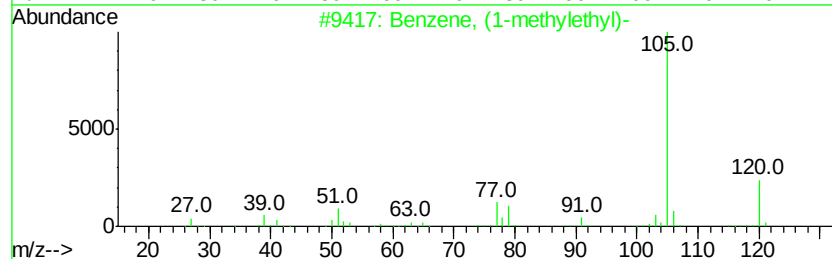
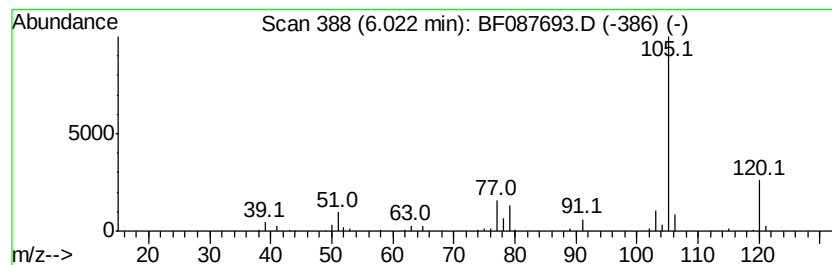
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, (1-methylethyl)- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.02	2.41 ng	91733	1,4-Dichlorobenzene-d4	6.86

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-2-methyl-	120	C9H12	000611-14-3	80
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	78
5		Mesitylene	120	C9H12	000108-67-8	78



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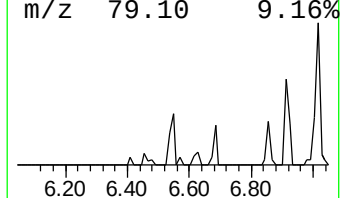
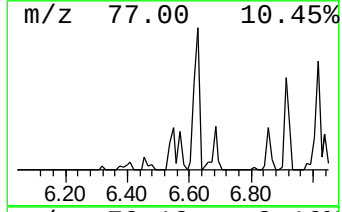
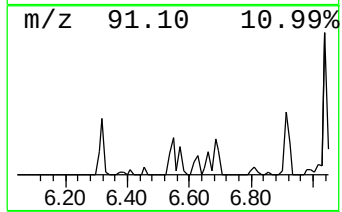
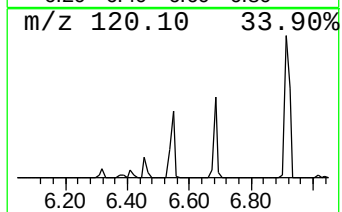
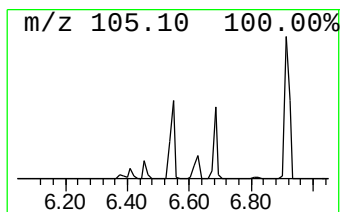
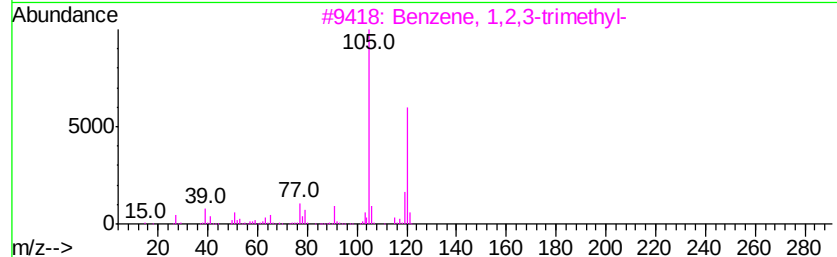
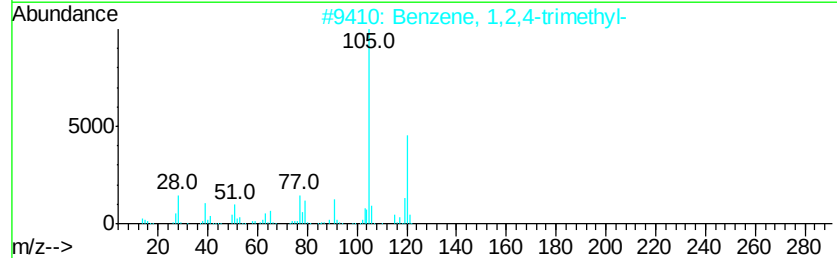
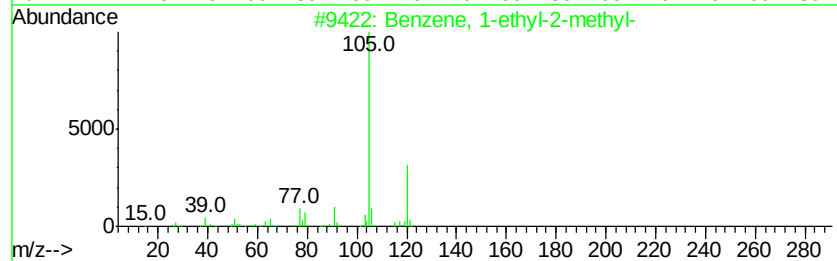
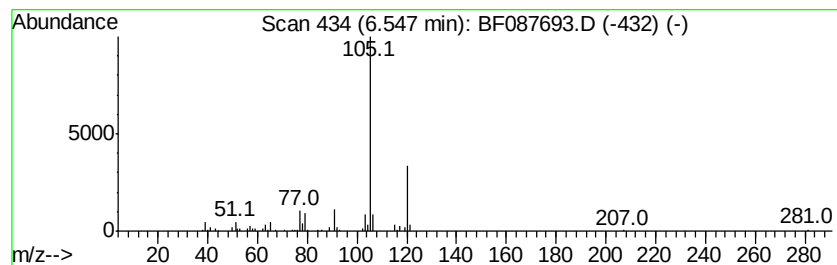
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	3.84 ng	146110	1,4-Dichlorobenzene-d4	6.86

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
Data File : BF087693.D
Acq On : 5 Jun 2016 7:11
Operator : UM/SJ
Sample : H3344-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

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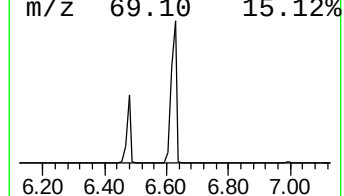
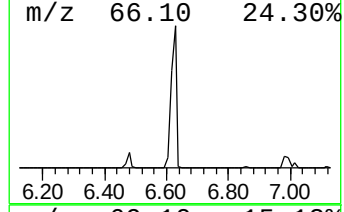
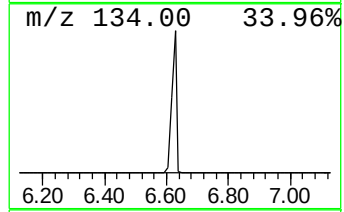
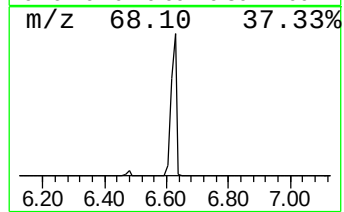
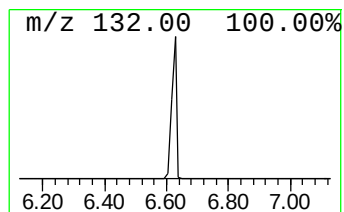
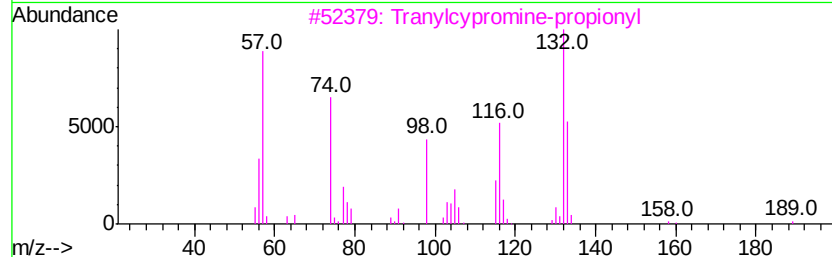
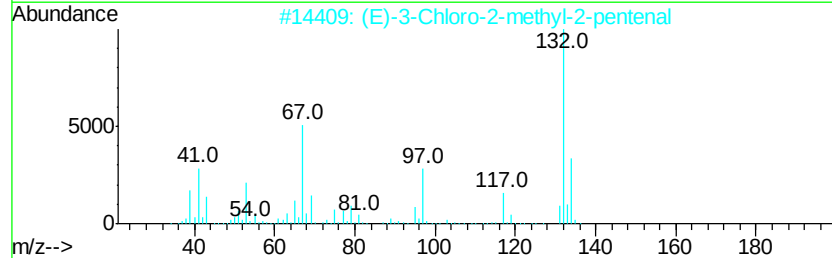
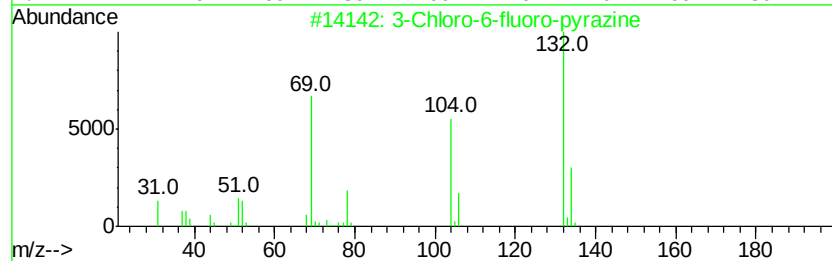
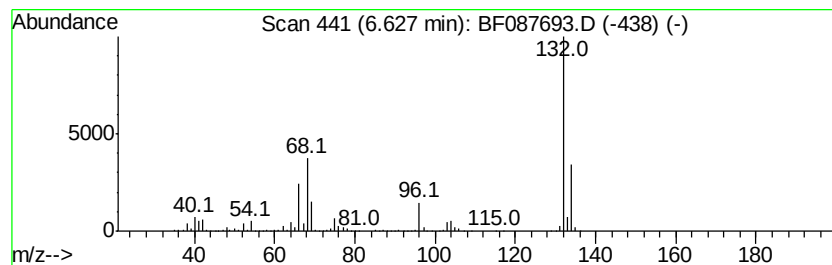
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	85.20 ng	3245000	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
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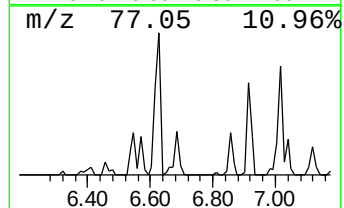
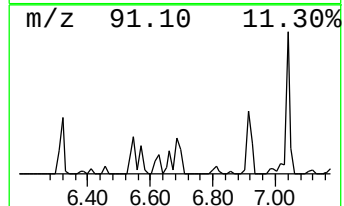
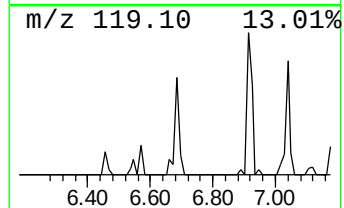
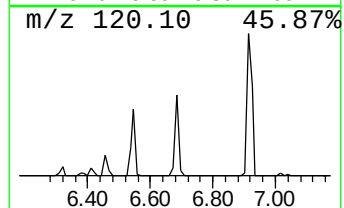
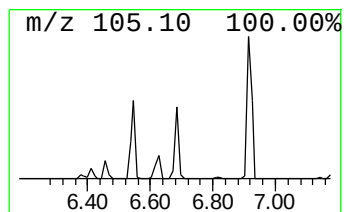
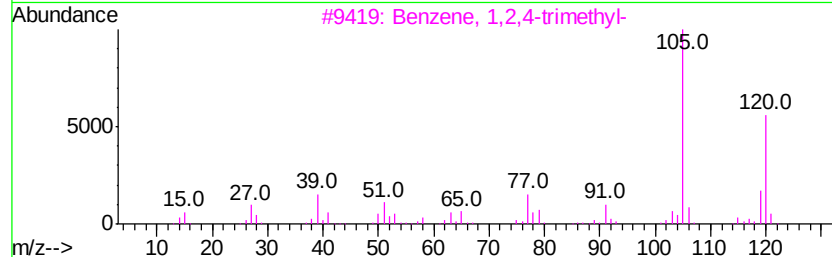
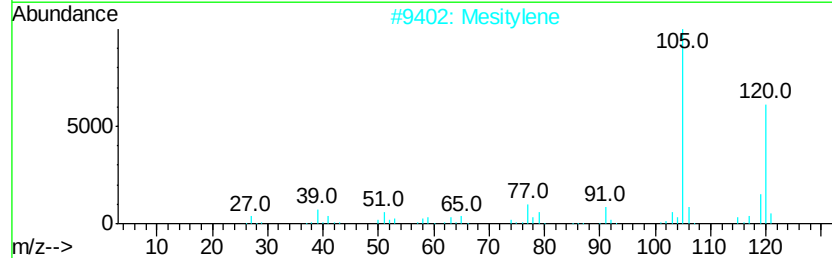
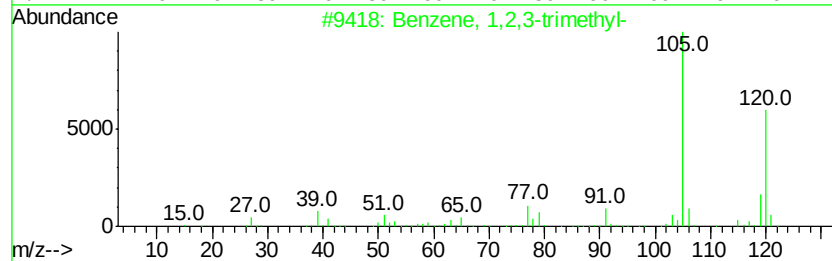
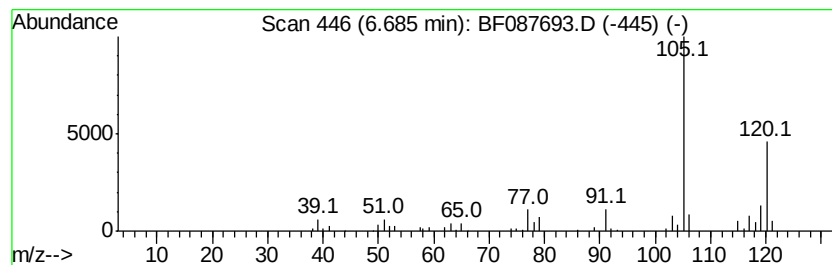
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	3.63 ng	138256	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Mesitylene	120	C9H12	000108-67-8	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
Data File : BF087693.D
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Misc :
ALS Vial : 35 Sample Multiplier: 1

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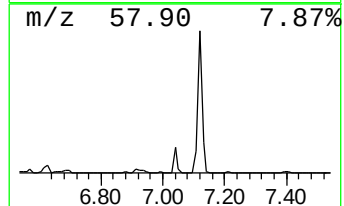
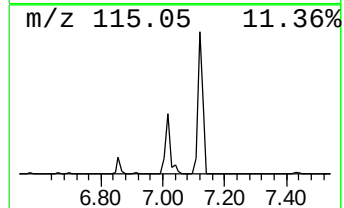
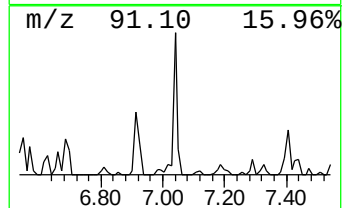
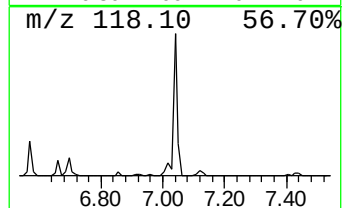
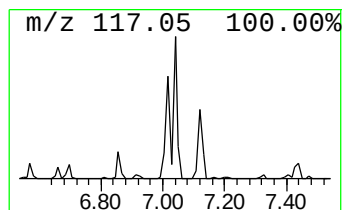
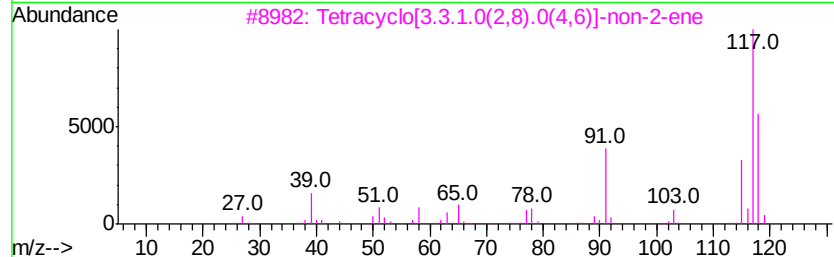
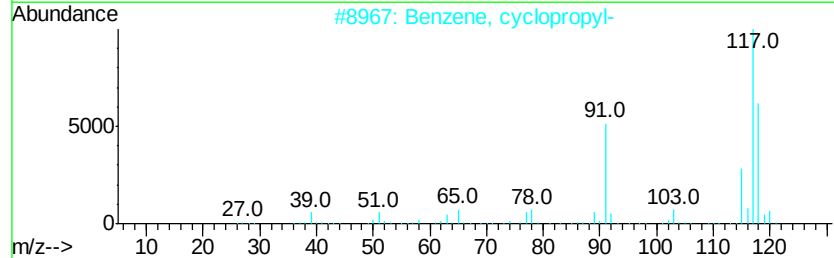
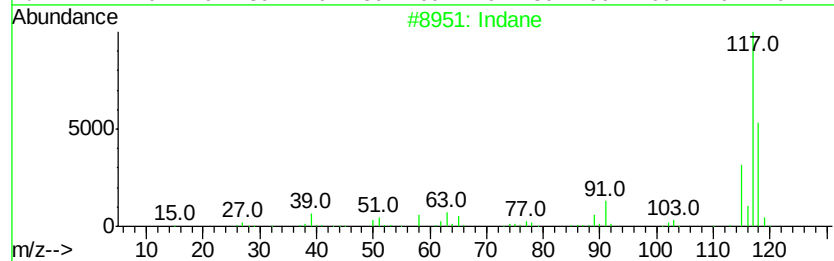
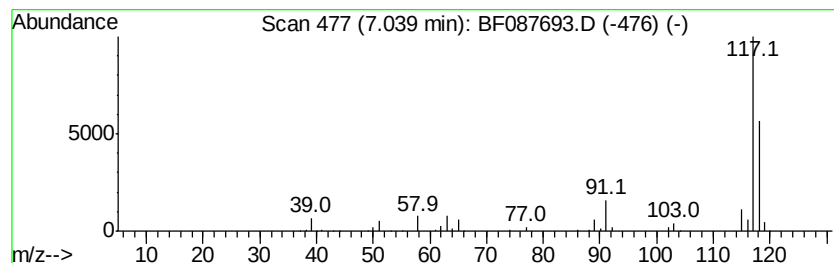
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	9.22 ng	351320	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	93
2		Benzene, cyclopropyl-	118	C9H10	000873-49-4	86
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	86
4		Benzene, 1-propenyl-	118	C9H10	000637-50-3	72
5		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
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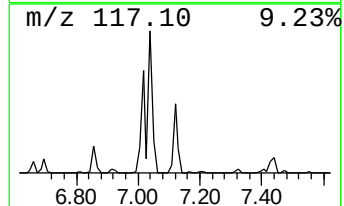
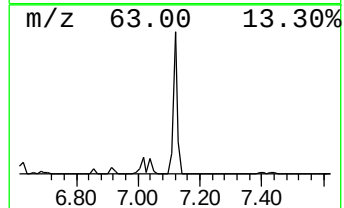
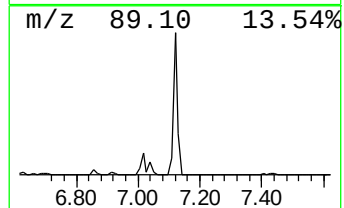
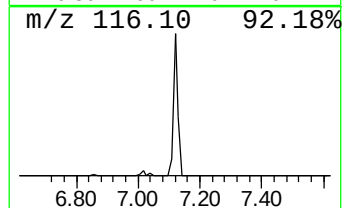
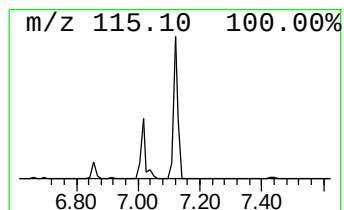
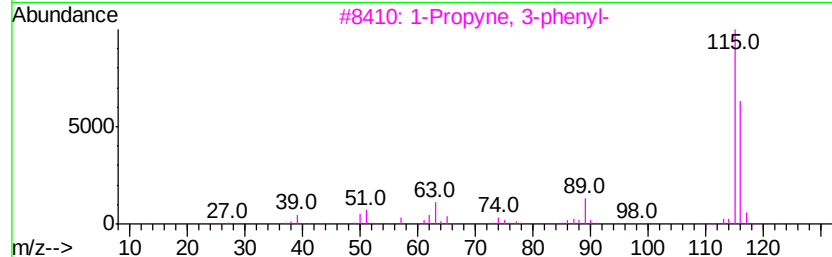
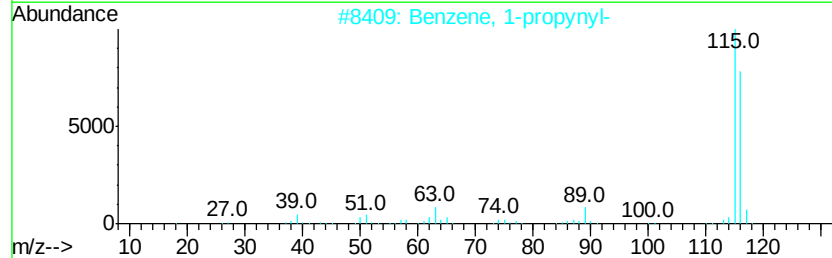
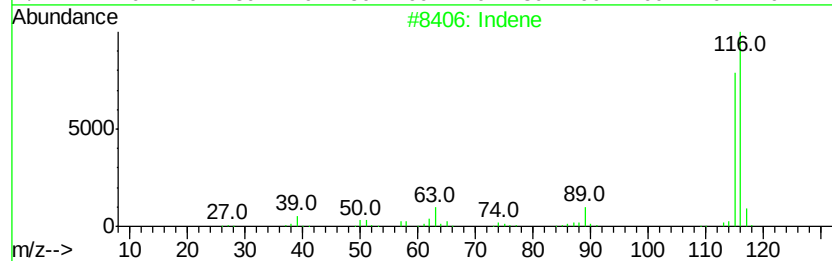
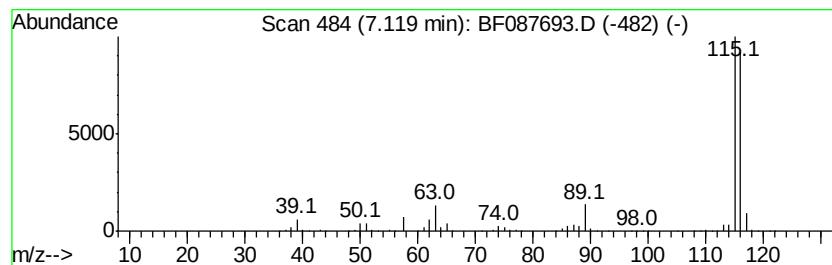
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Indene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	61.48 ng	2341640	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	97
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3		1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
4		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
5		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
Data File : BF087693.D
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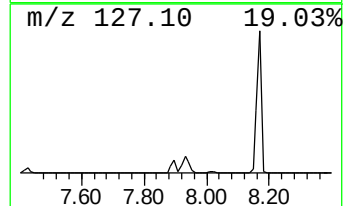
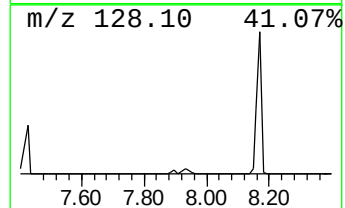
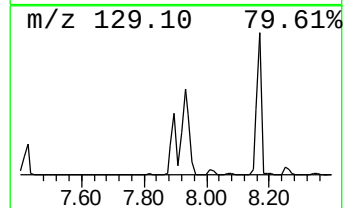
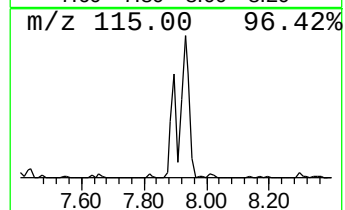
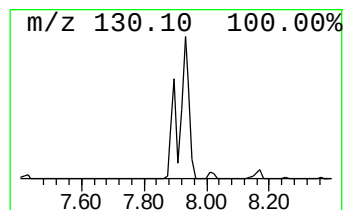
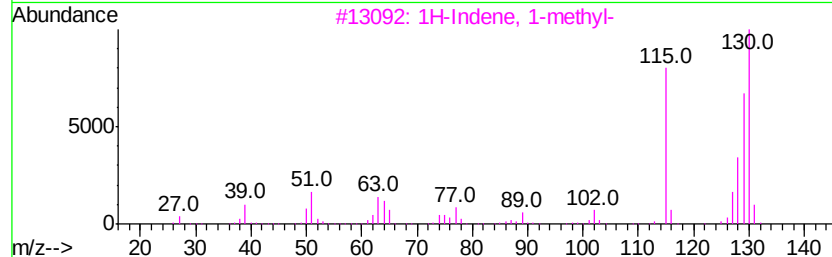
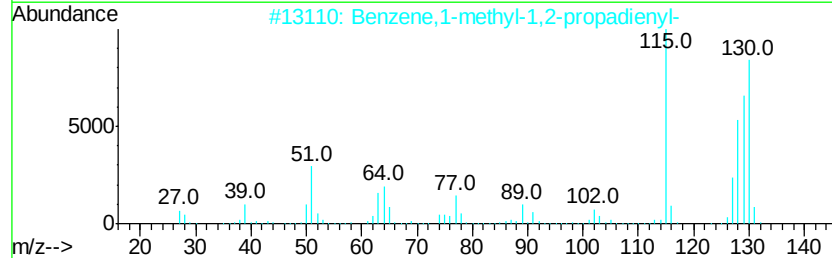
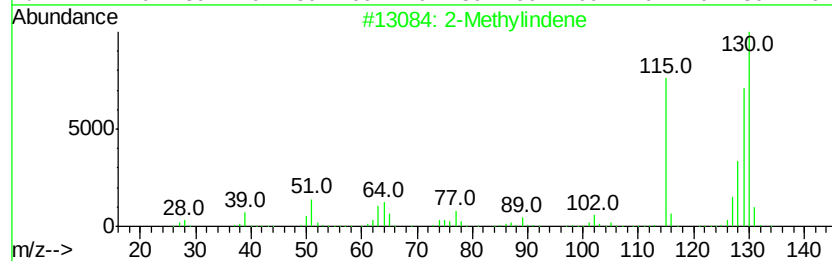
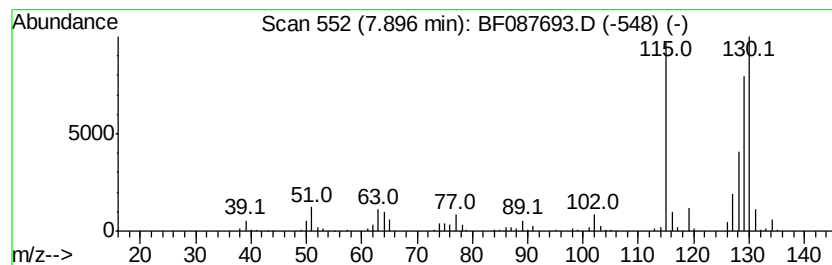
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 2-Methylindene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	3.07 ng	488958	Naphthalene-d8	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methylindene	130	C10H10	002177-47-1	96
2		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
3		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	94
4		Benzene, 1-butynyl-	130	C10H10	000622-76-4	94
5		1H-Indene, 3-methyl-	130	C10H10	000767-60-2	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
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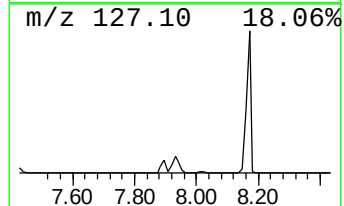
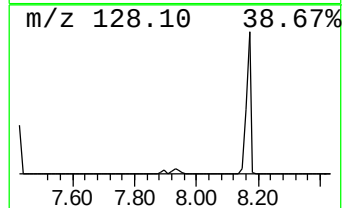
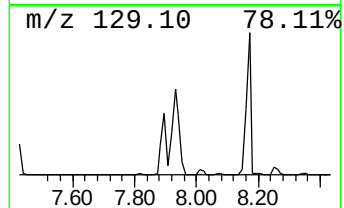
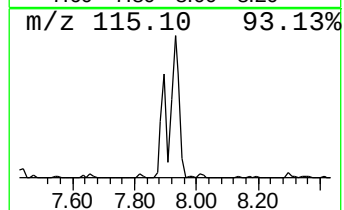
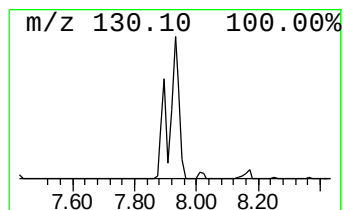
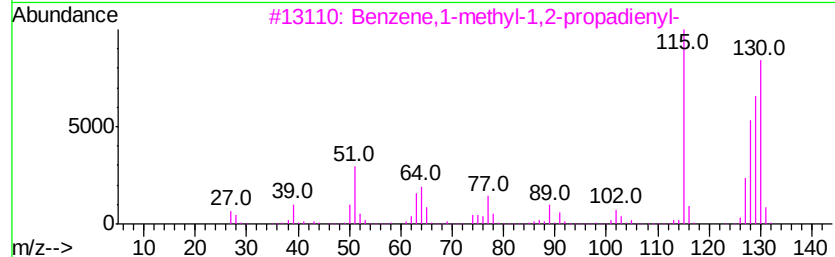
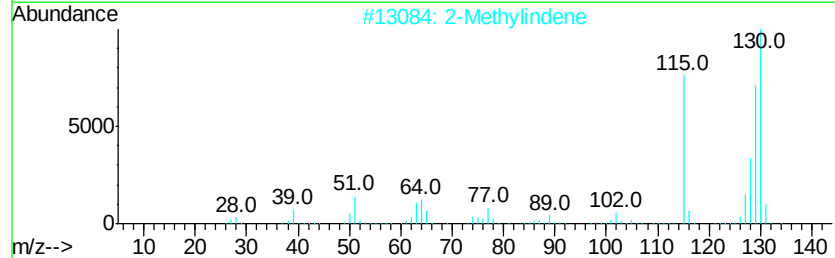
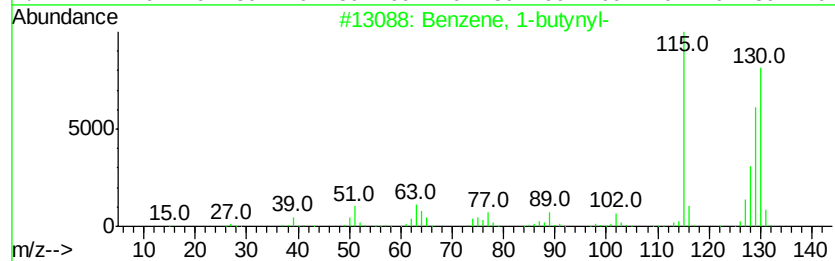
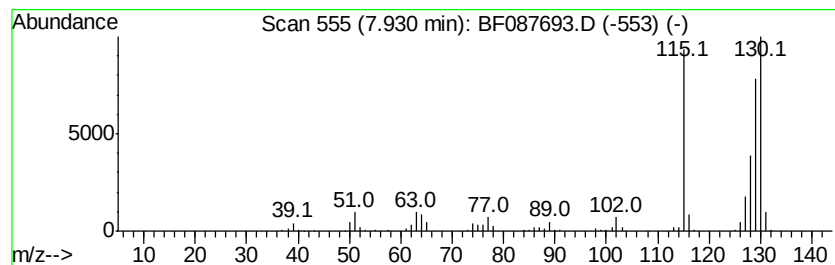
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TIC Integration Parameters: LSCINT.P

Peak Number 17 Benzene, 1-butynyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	4.38 ng	696379	Napthalene-d8	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-butynyl-	130	C10H10	000622-76-4	96
2		2-Methylindene	130	C10H10	002177-47-1	95
3		Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
4		Cycloprop[alindene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93
5		1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
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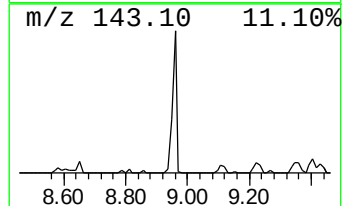
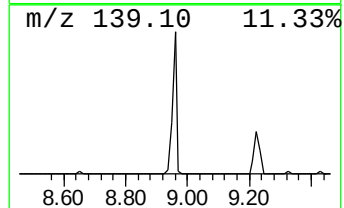
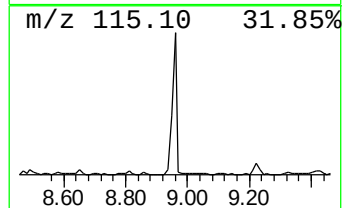
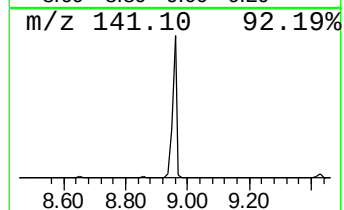
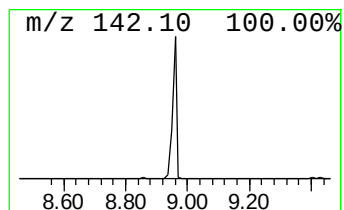
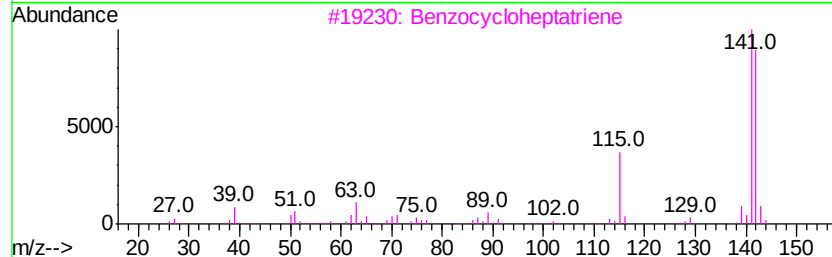
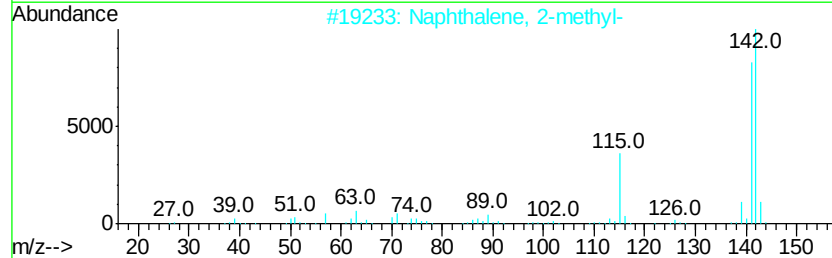
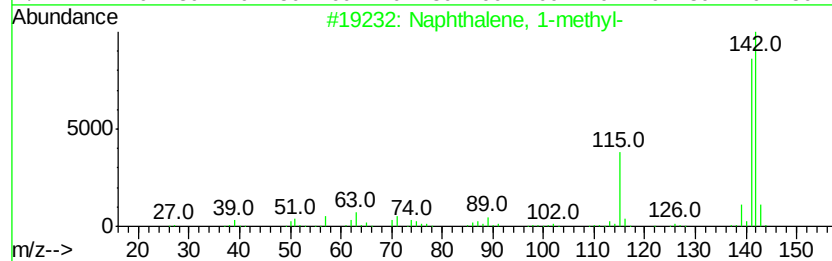
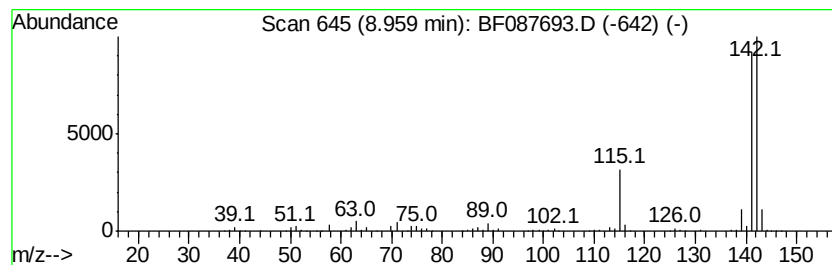
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ClientSampleId :
MW-18

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Naphthalene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.96	4.65 ng	740337	Naphthalene-d8	8.15	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
Data File : BF087693.D
Acq On : 5 Jun 2016 7:11
Operator : UM/SJ
Sample : H3344-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-18

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	2.15	6.6	ng	250824	1	6.86	761740	20.0
2-Pentanone, 4-hy...	5.04	4.8	ng	181906	1	6.86	761740	20.0
Benzene, 1,3-dime...	5.69	18.2	ng	694681	1	6.86	761740	20.0
Benzene, (1-methy...	6.02	2.4	ng	91733	1	6.86	761740	20.0
Benzene, 1-ethyl-...	6.55	3.8	ng	146110	1	6.86	761740	20.0
unknown6.63	6.63	85.2	ng	3245000	1	6.86	761740	20.0
Benzene, 1,2,3-tr...	6.68	3.6	ng	138256	1	6.86	761740	20.0
Indane	7.04	9.2	ng	351320	1	6.86	761740	20.0
Indene	7.12	61.5	ng	2341640	1	6.86	761740	20.0
2-Methylindene	7.90	3.1	ng	488958	2	8.15	3182210	20.0
Benzene, 1-butynyl-	7.93	4.4	ng	696379	2	8.15	3182210	20.0
Naphthalene, 1-me...	8.96	4.7	ng	740337	2	8.15	3182210	20.0