

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
 Data File : BF081676.D
 Acq On : 17 Sep 2015 13:17
 Operator : UM/IZ
 Sample : G3669-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-91415-WATER

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-BF090115.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	1.224	9	12	15	rBV	2734694	2953073	17.23%	5.228%
2	1.532	37	39	57	rBV	155736	480218	2.80%	0.850%
3	2.504	120	124	130	rVB	475898	792912	4.63%	1.404%
4	4.424	289	292	307	rBV	152201	317731	1.85%	0.562%
5	4.904	329	334	337	rBV	621358	1279679	7.47%	2.265%
6	5.121	349	353	355	rBV	414213	872679	5.09%	1.545%
7	5.327	368	371	380	rVB	88427	200323	1.17%	0.355%
8	5.647	394	399	407	rVB	594685	1172364	6.84%	2.075%
9	7.144	527	530	533	rBV	128441	208768	1.22%	0.370%
10	7.602	565	570	575	rBV2	55129	176075	1.03%	0.312%
11	7.682	575	577	582	rVV	199606	383305	2.24%	0.679%
12	7.830	585	590	594	rVB2	239406	731853	4.27%	1.296%
13	8.173	617	620	628	rVB	161814	283716	1.66%	0.502%
14	8.505	645	649	651	rVV	369131	747220	4.36%	1.323%
15	8.562	651	654	658	rVB	201781	383901	2.24%	0.680%
16	8.779	665	673	676	rBV	2939847	9270038	54.09%	16.410%
17	9.476	730	734	745	rVB	351803	722939	4.22%	1.280%
18	10.562	821	829	831	rBV	246674	659220	3.85%	1.167%
19	10.619	831	834	835	rVV	271322	541214	3.16%	0.958%
20	10.653	835	837	841	rVV	178530	297960	1.74%	0.527%
21	11.202	869	885	887	rBV	4043506	17138842	100.00%	30.339%
22	11.271	887	891	893	rVB	912698	1481805	8.65%	2.623%
23	11.556	913	916	927	rBV	215228	396530	2.31%	0.702%
24	12.642	1008	1011	1018	rVB	112931	206702	1.21%	0.366%
25	12.791	1018	1024	1027	rBV	1369501	2993170	17.46%	5.298%
26	13.008	1035	1043	1046	rVB	1258415	2903566	16.94%	5.140%
27	13.717	1100	1105	1108	rBV	658989	1096372	6.40%	1.941%
28	13.911	1119	1122	1126	rBV	162302	272403	1.59%	0.482%
29	14.117	1136	1140	1145	rVB	93730	187937	1.10%	0.333%
30	14.265	1147	1153	1158	rBV3	118522	377751	2.20%	0.669%
31	14.448	1165	1169	1172	rBV	224792	430315	2.51%	0.762%
32	14.505	1172	1174	1177	rVB	141003	207570	1.21%	0.367%
33	14.711	1186	1192	1195	rBV	78165	216431	1.26%	0.383%
34	14.860	1200	1205	1207	rBV	740134	1331748	7.77%	2.357%

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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.202	1230	1235	1238	rBV	218678	383224	2.24%	0.678%
36	15.271	1238	1241	1245	rVV	245283	386613	2.26%	0.684%
37	16.425	1337	1342	1346	rBV2	79069	256597	1.50%	0.454%
38	16.517	1346	1350	1355	rVB	331881	658572	3.84%	1.166%
39	17.100	1397	1401	1407	rVB2	159593	300740	1.75%	0.532%
40	18.700	1537	1541	1543	rBV	222721	399797	2.33%	0.708%
41	18.757	1543	1546	1549	rVB	446262	790492	4.61%	1.399%
42	22.072	1833	1836	1838	rBV	801174	874014	5.10%	1.547%
43	23.352	1942	1948	1950	rBV	273108	491462	2.87%	0.870%
44	24.781	2071	2073	2076	rBV	216576	233267	1.36%	0.413%

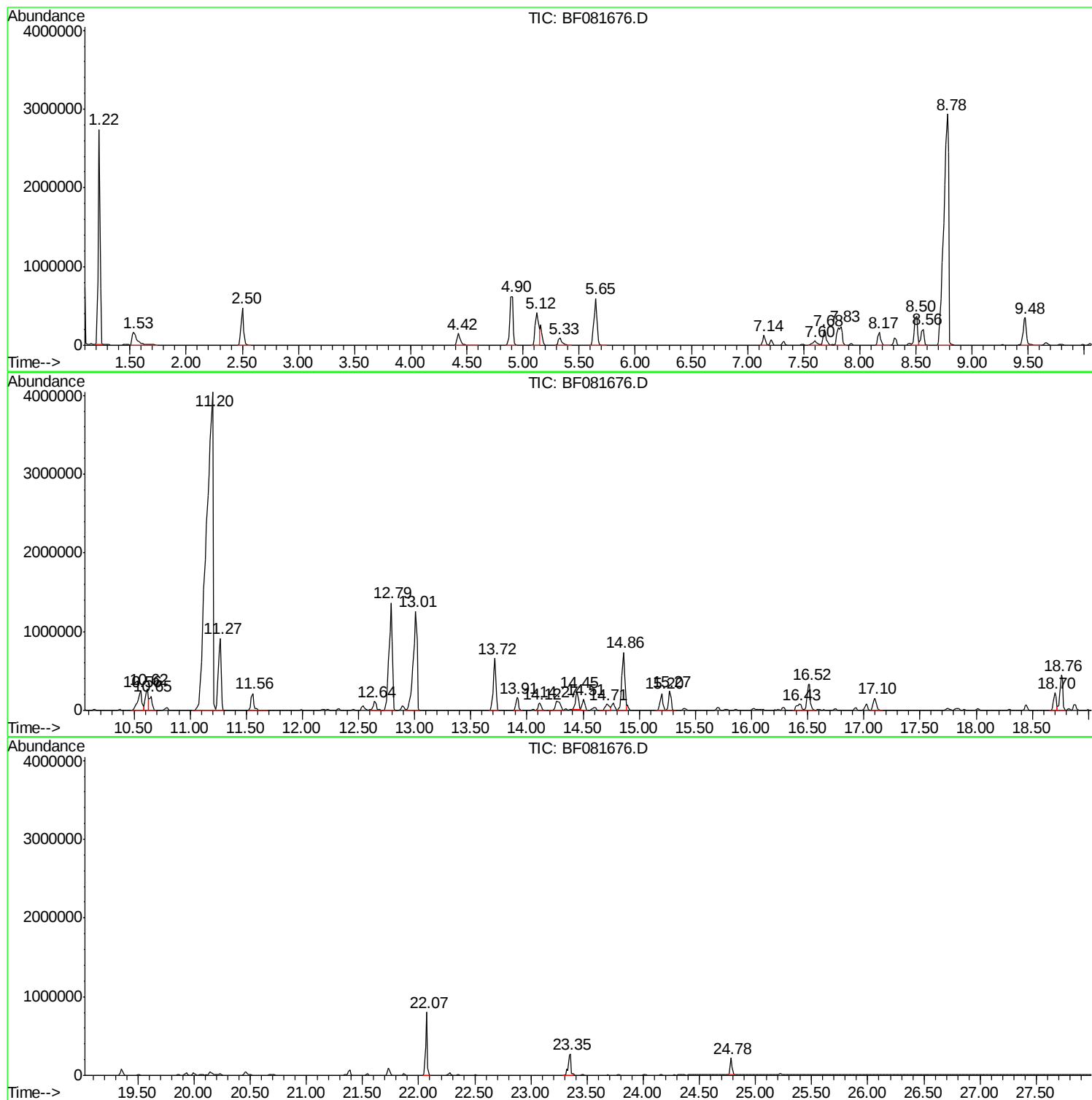
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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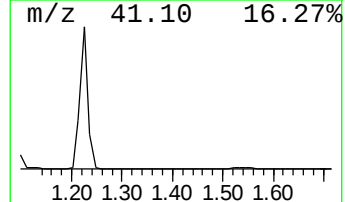
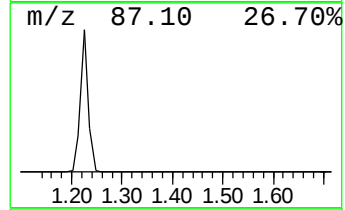
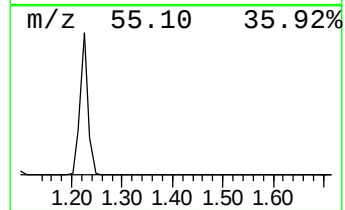
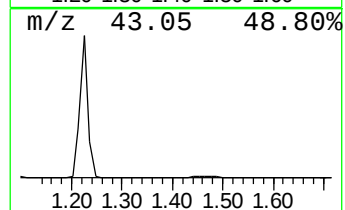
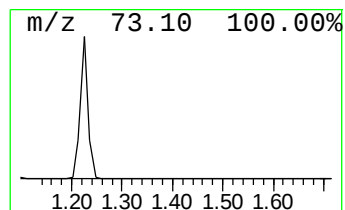
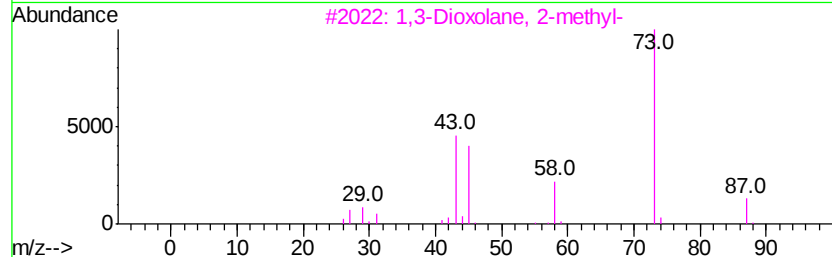
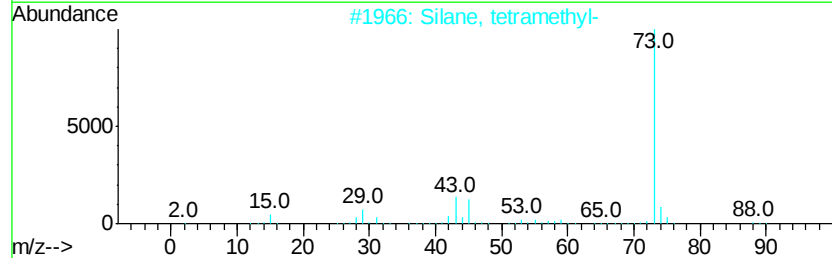
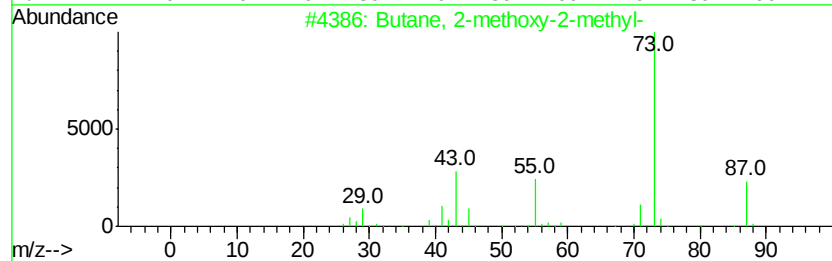
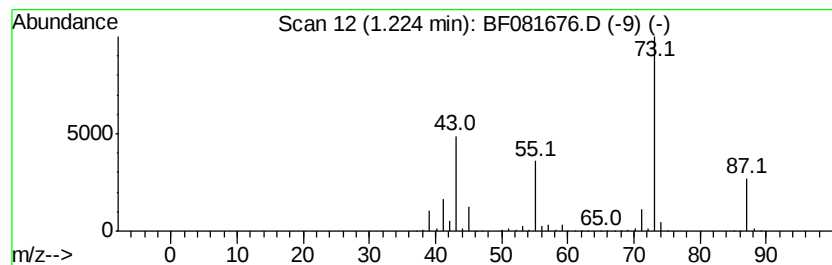
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.22	208.17 ng	2953070	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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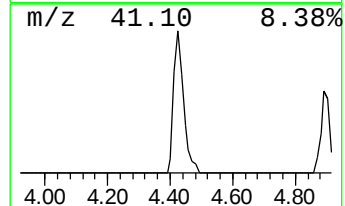
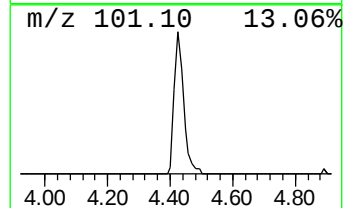
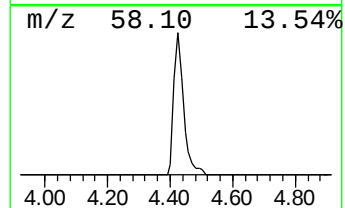
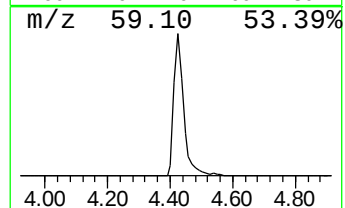
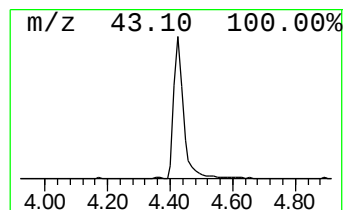
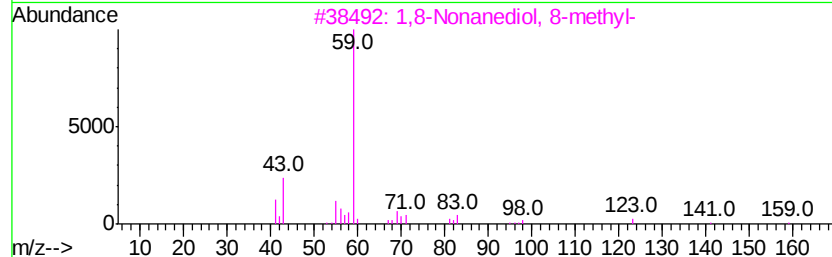
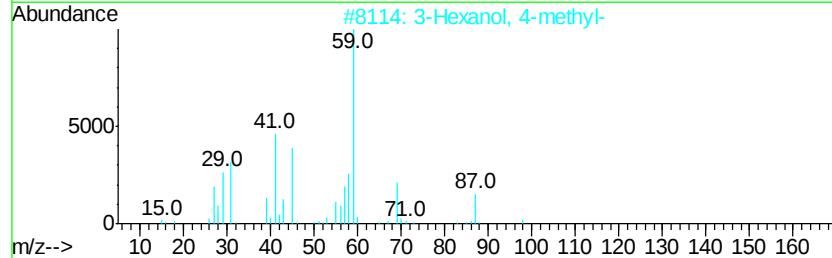
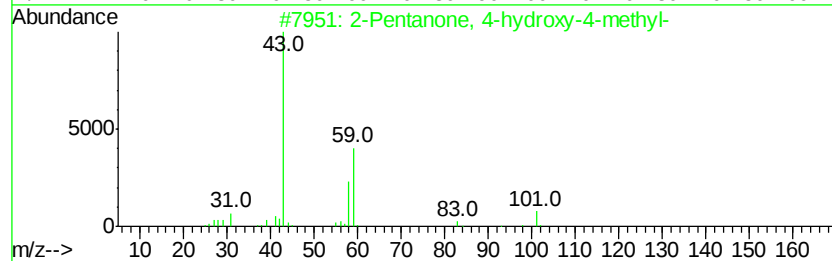
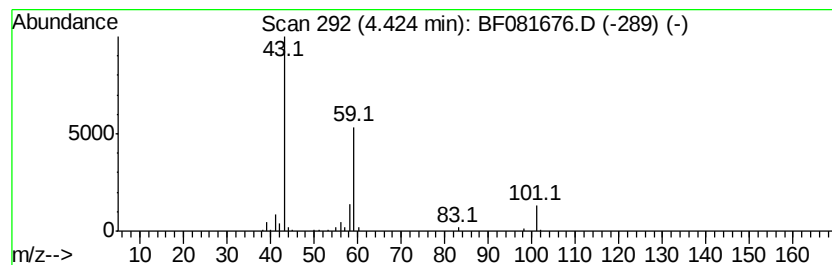
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.42	22.40 ng	317731	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9
4		3-Octanol	130	C8H18O	000589-98-0	9
5		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	9



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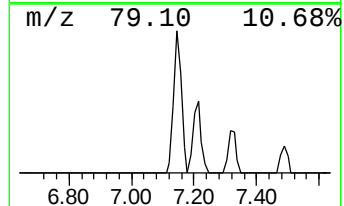
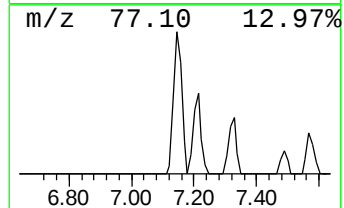
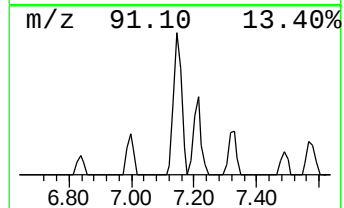
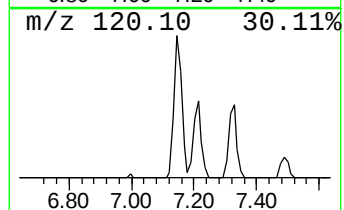
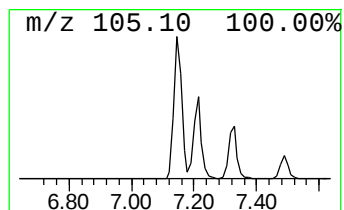
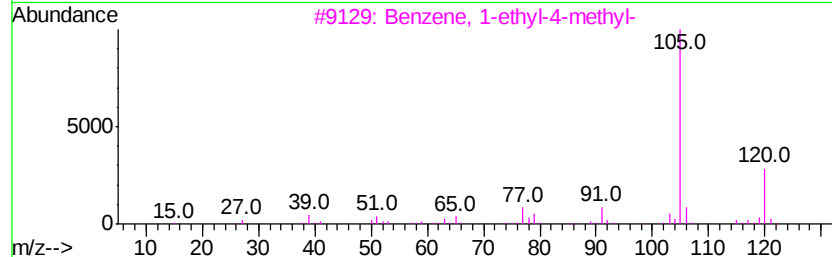
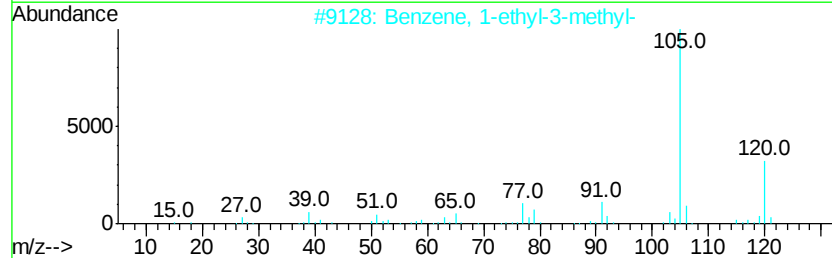
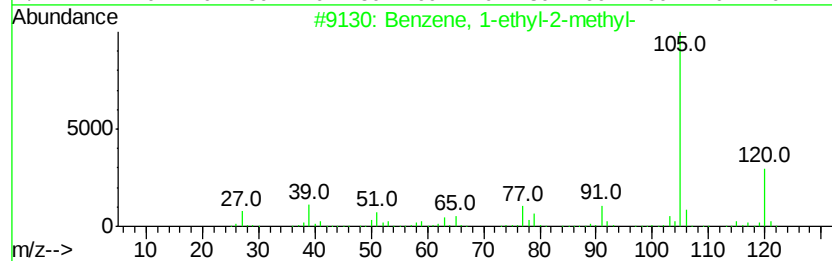
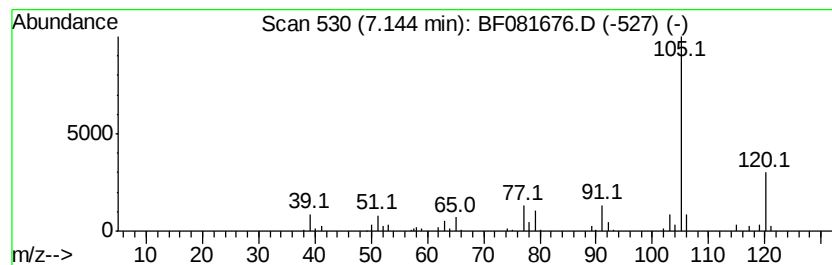
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.14	14.72 ng	208768	1,4-Dichlorobenzene-d4	8.17

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



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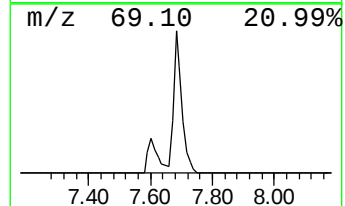
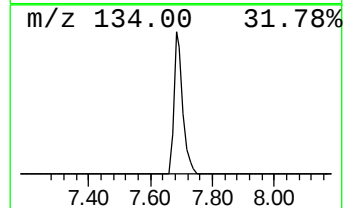
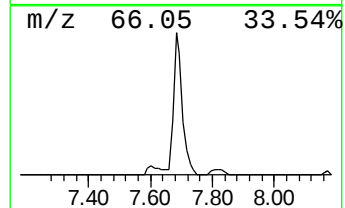
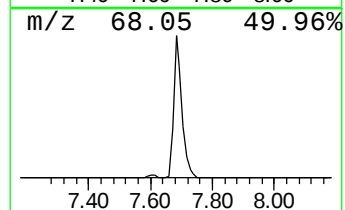
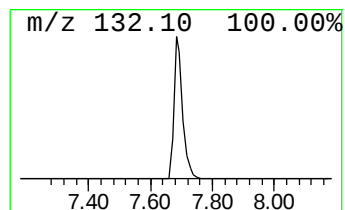
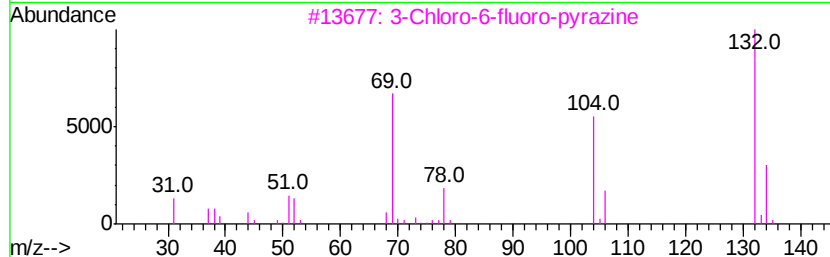
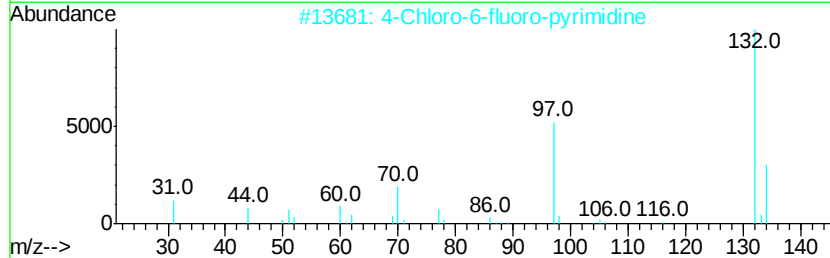
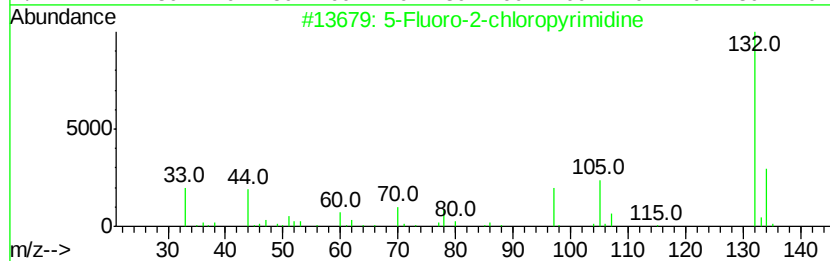
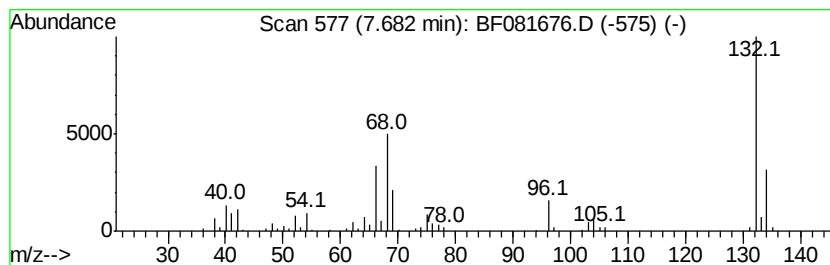
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown7.68 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	27.02 ng	383305	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	25
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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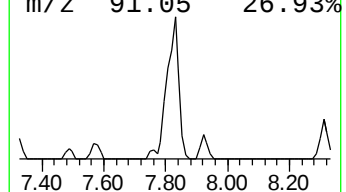
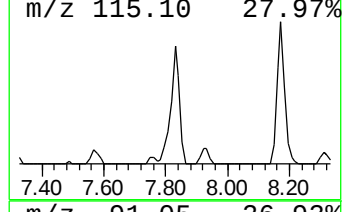
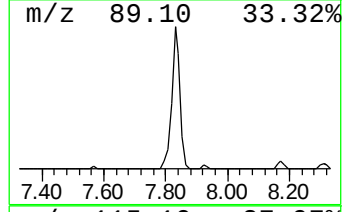
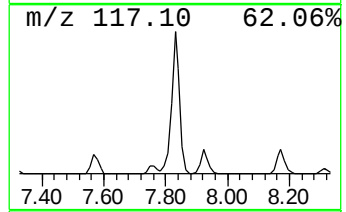
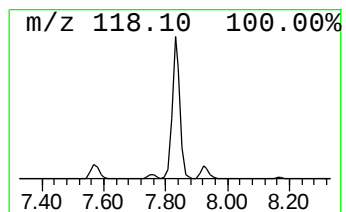
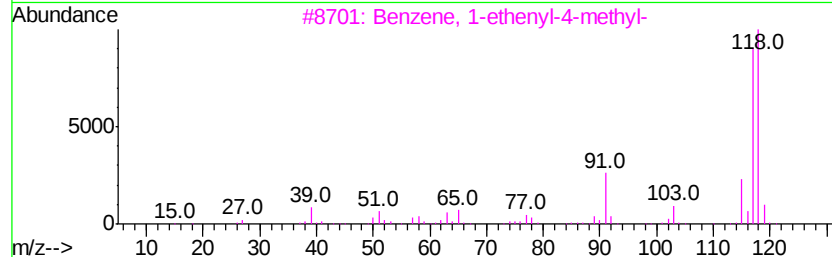
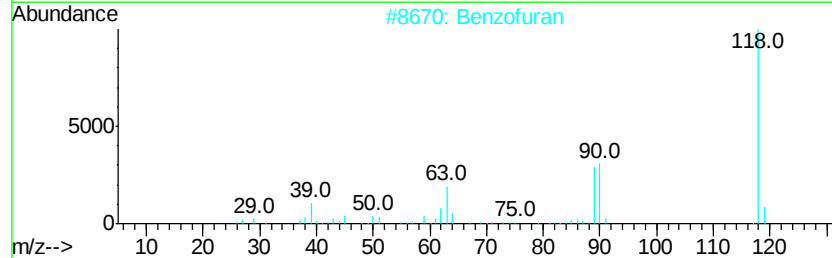
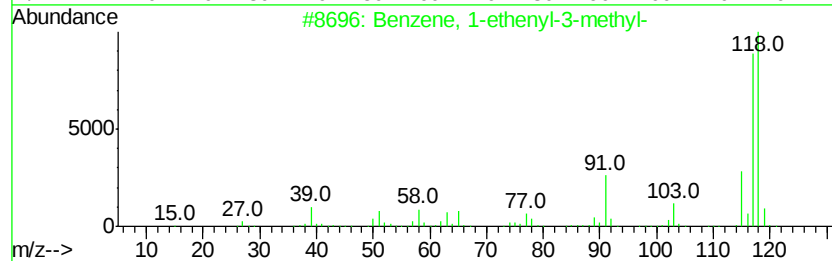
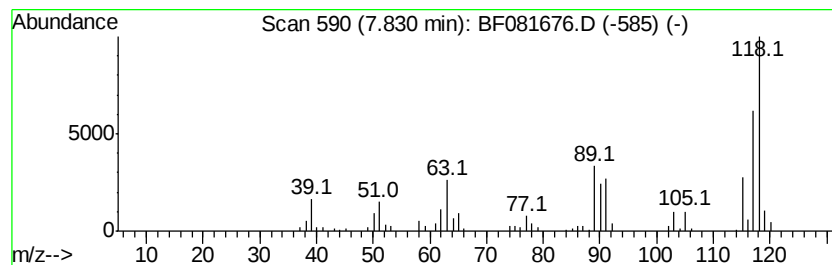
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1-ethenyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	51.59 ng	731853	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	87
2		Benzofuran	118	C8H6O	000271-89-6	62
3		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	55
4		3-Pyridineacetonitrile	118	C7H6N2	006443-85-2	53
5		Indane	118	C9H10	000496-11-7	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
Data File : BF081676.D
Acq On : 17 Sep 2015 13:17
Operator : UM/IZ
Sample : G3669-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-91415-WATER

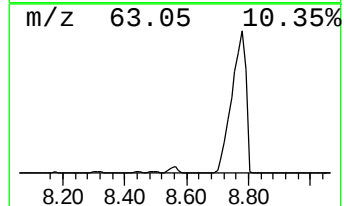
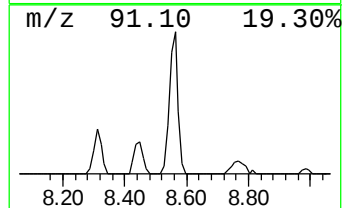
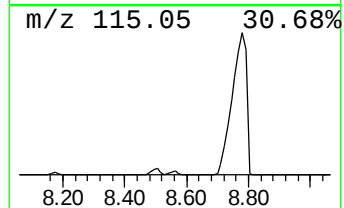
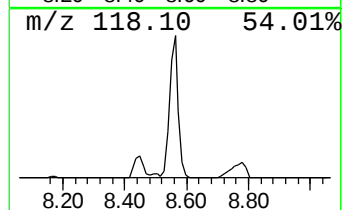
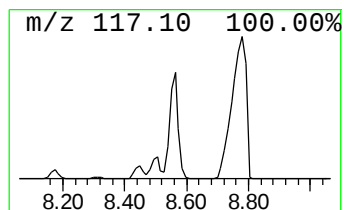
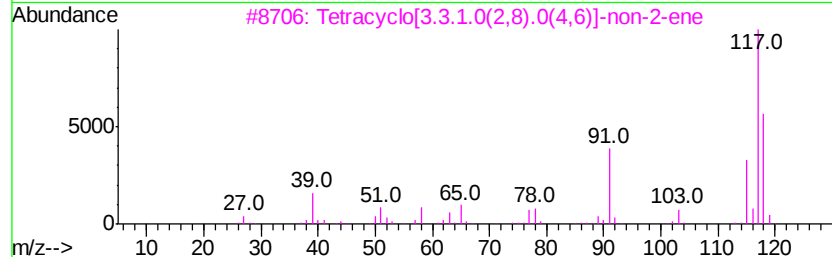
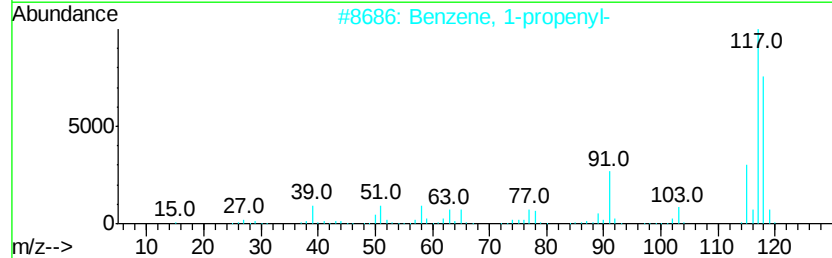
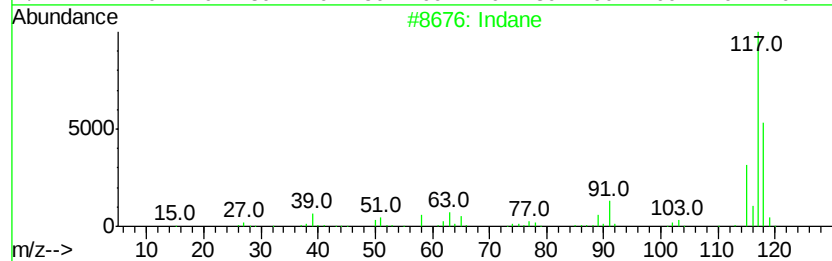
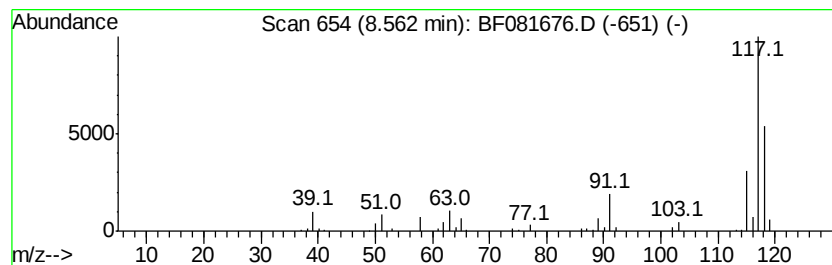
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	27.06 ng	383901	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	91
2		Benzene, 1-propenyl-	118	C9H10	000637-50-3	76
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
4		Deltacyclene	118	C9H10	007785-10-6	68
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	60



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
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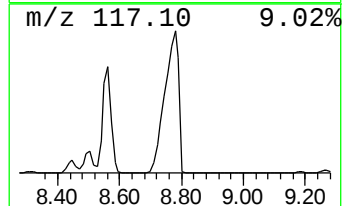
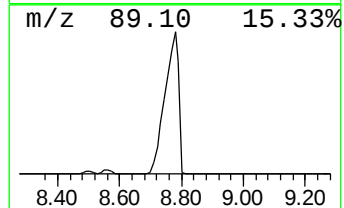
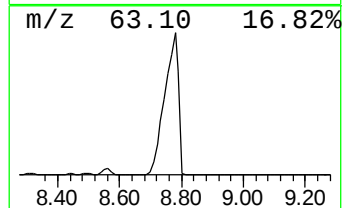
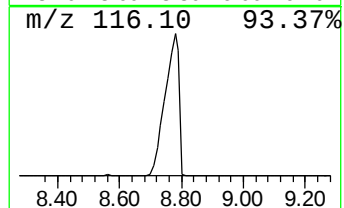
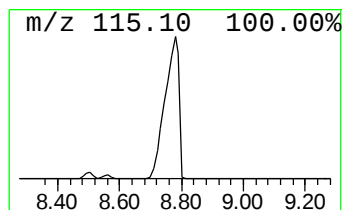
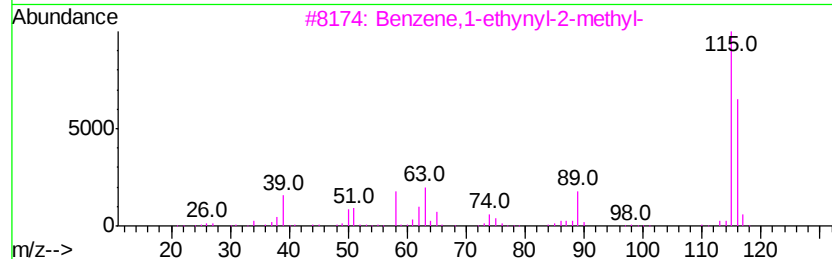
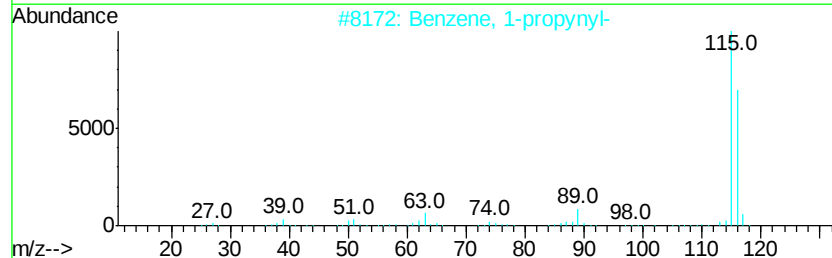
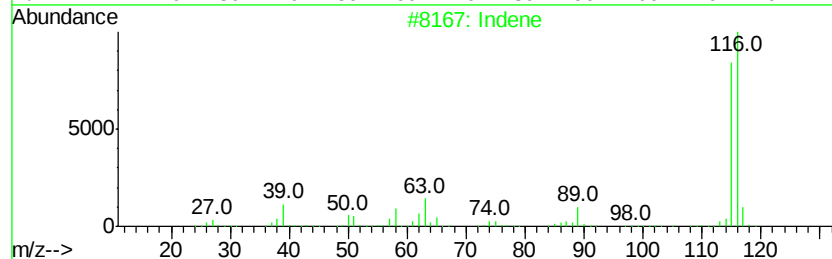
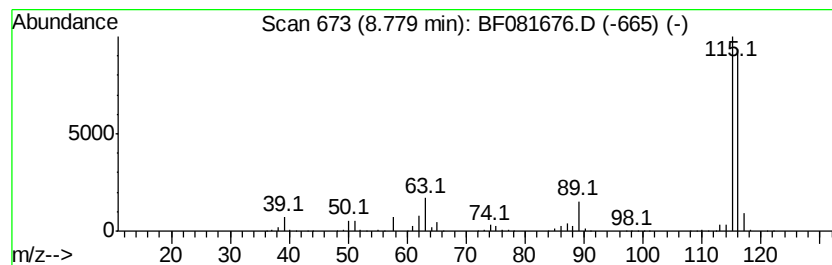
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.78	653.47 ng	9270040	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	95
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
3		Benzene,1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91
4		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
5		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
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Operator : UM/IZ
Sample : G3669-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

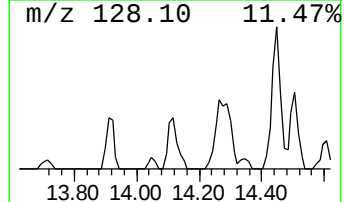
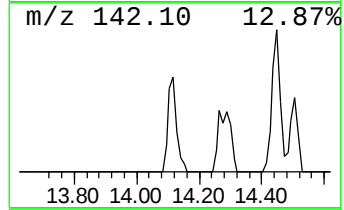
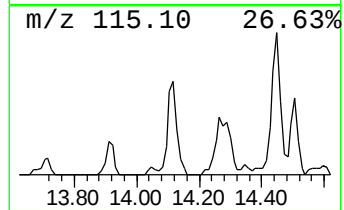
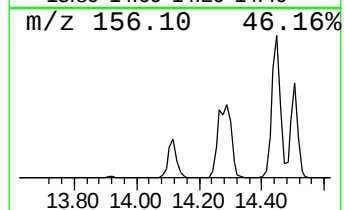
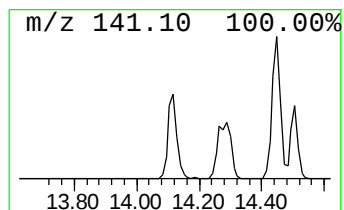
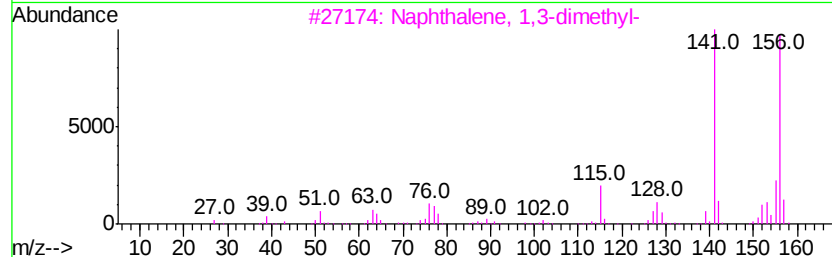
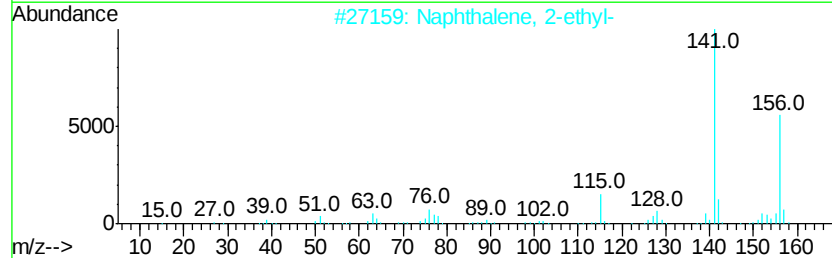
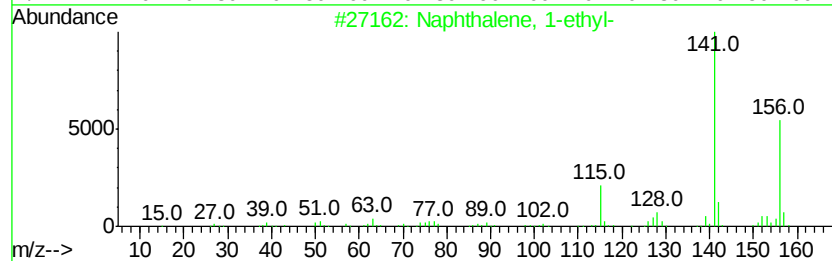
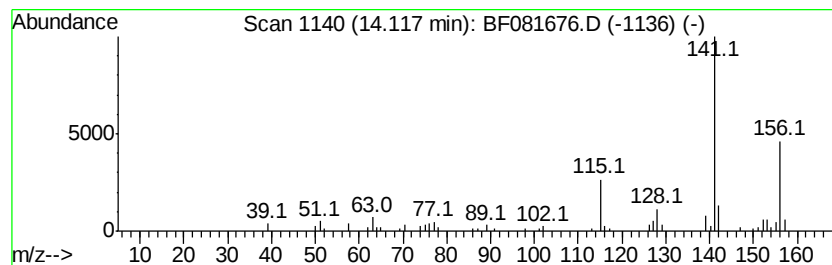
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 1-ethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	9.81 ng	187937	Acenaphthene-d10	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-ethyl-	156 C12H12	001127-76-0 95
2		Naphthalene, 2-ethyl-	156 C12H12	000939-27-5 94
3		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 91
4		2-Thiophenecarboxylic acid, 5-et...	156 C7H8O2S	023229-72-3 59
5		3',4'-Difluoroacetophenone	156 C8H6F2O	000369-33-5 53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
Data File : BF081676.D
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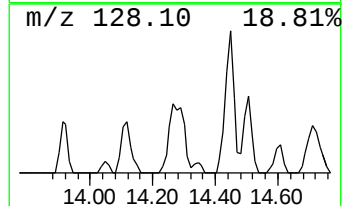
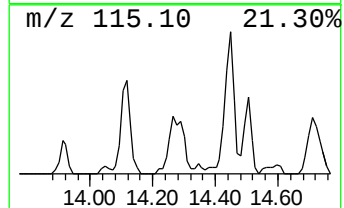
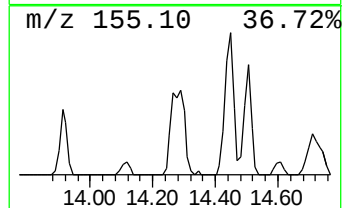
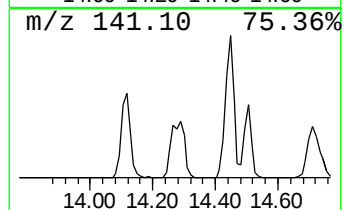
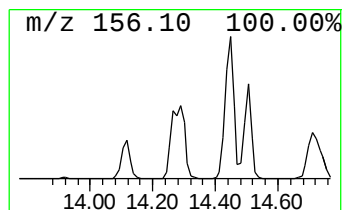
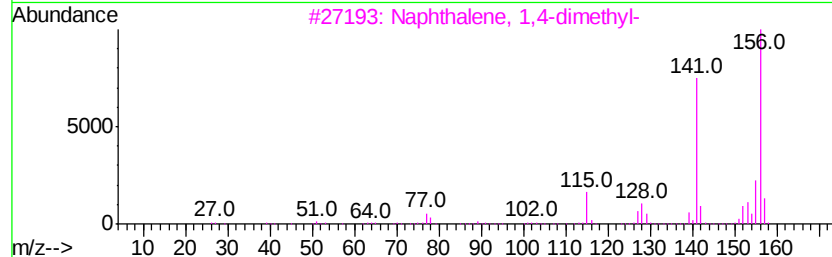
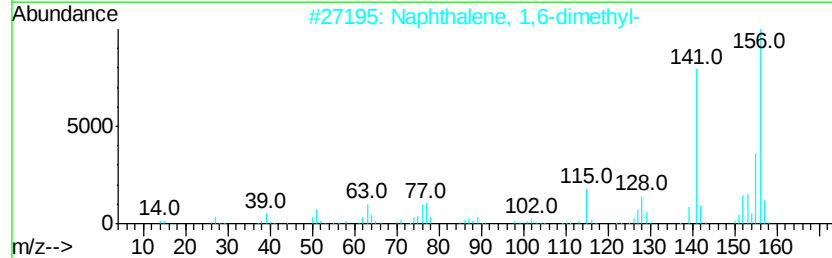
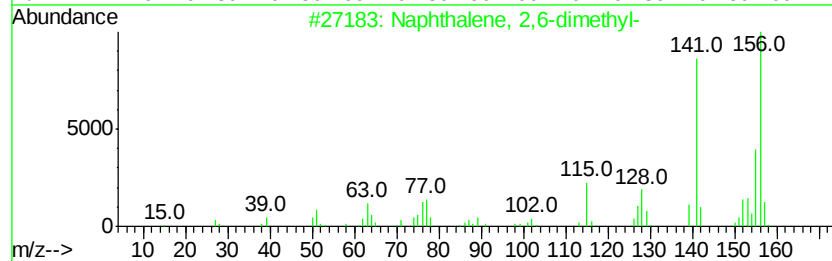
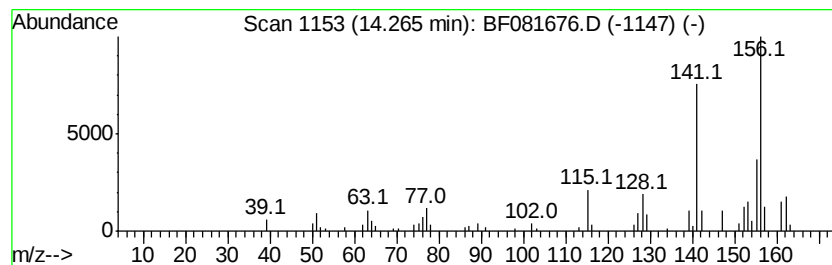
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WC-91415-WATER

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 2,6-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	19.71 ng	377751	Acenaphthene-d10	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 96
2		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 95
3		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 95
4		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 95
5		Naphthalene, 1,8-dimethyl-	156 C12H12	000569-41-5 93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
Data File : BF081676.D
Acq On : 17 Sep 2015 13:17
Operator : UM/IZ
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WC-91415-WATER

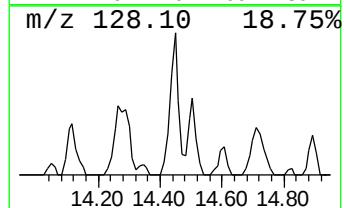
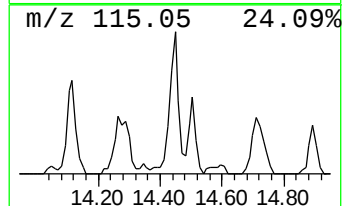
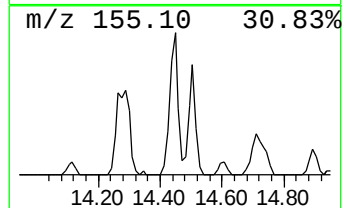
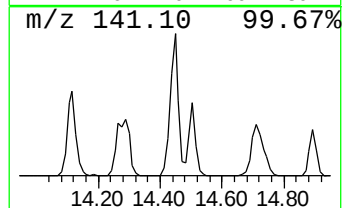
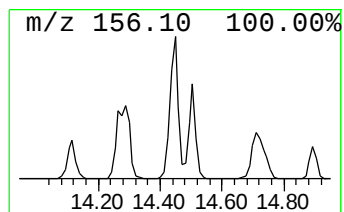
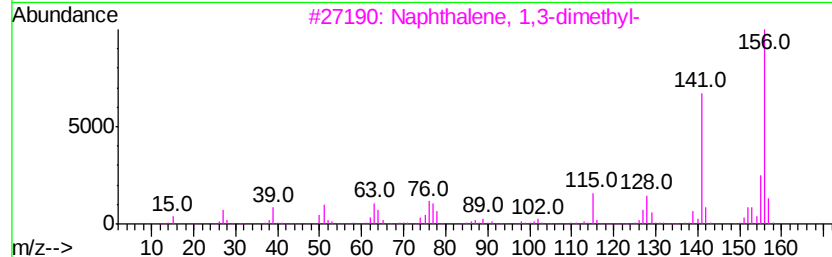
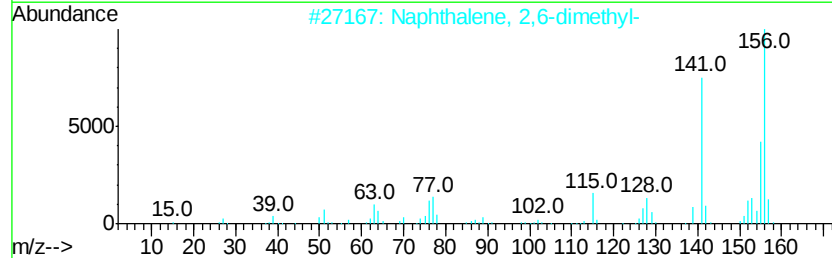
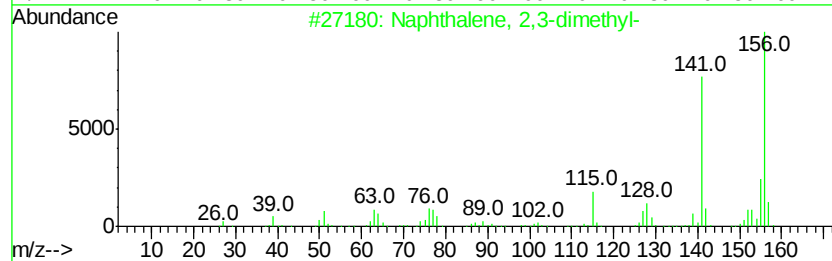
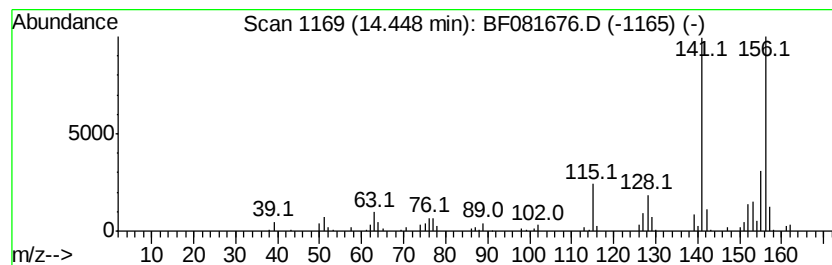
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	22.46 ng	430315	Acenaphthene-d10	15.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
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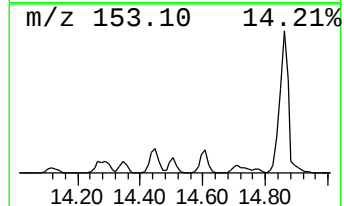
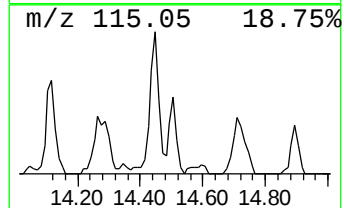
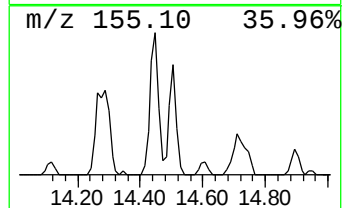
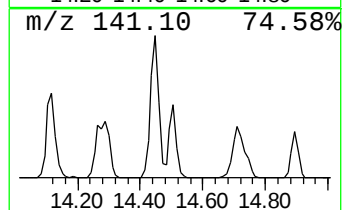
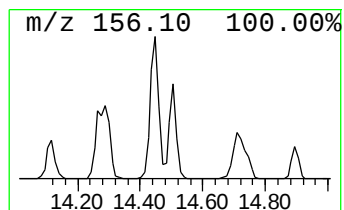
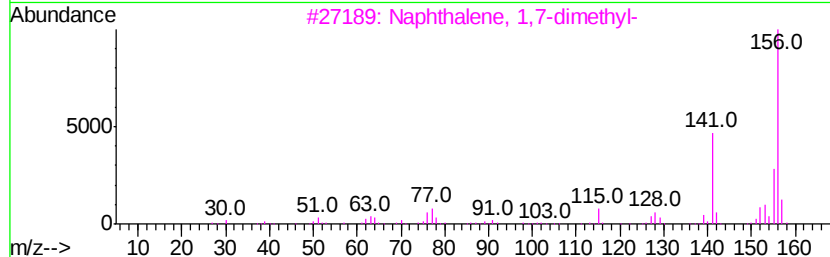
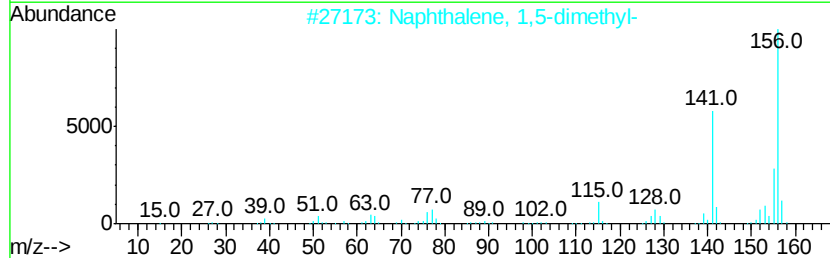
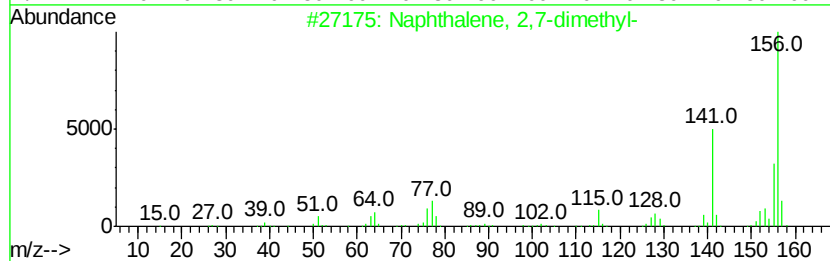
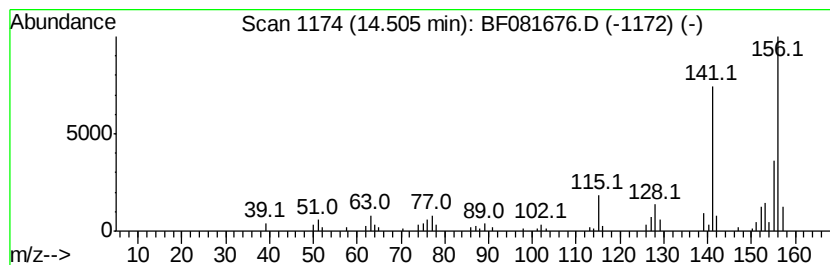
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Naphthalene, 2,7-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.51	10.83 ng	207570	Acenaphthene-d10	15.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
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Acq On : 17 Sep 2015 13:17
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Misc :
ALS Vial : 3 Sample Multiplier: 1

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ClientSampleId :
WC-91415-WATER

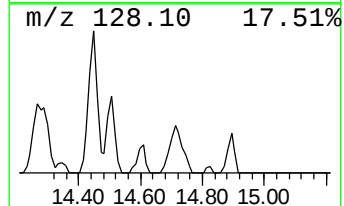
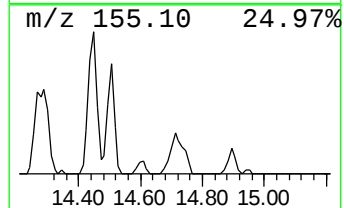
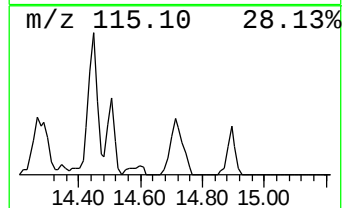
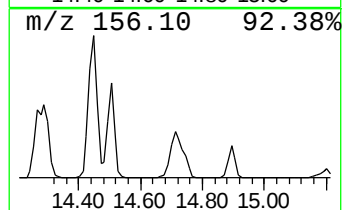
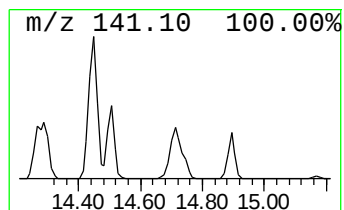
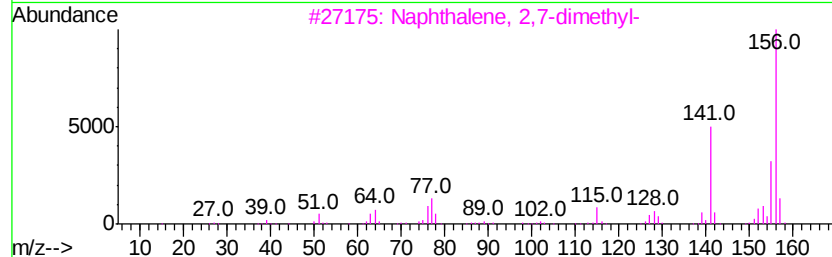
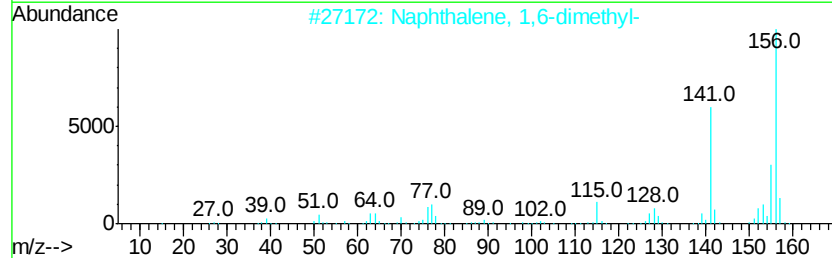
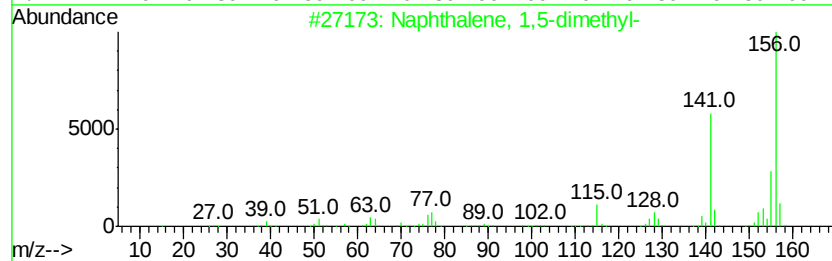
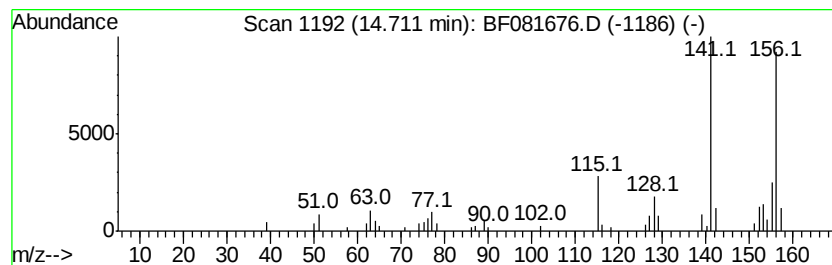
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Naphthalene, 1,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	11.30 ng	216431	Acenaphthene-d10	15.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
Data File : BF081676.D
Acq On : 17 Sep 2015 13:17
Operator : UM/IZ
Sample : G3669-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-91415-WATER

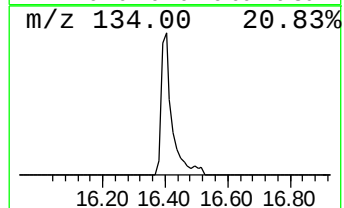
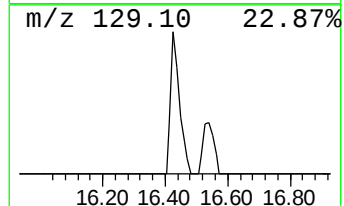
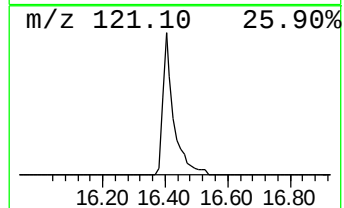
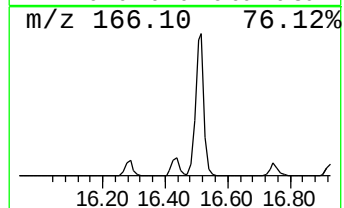
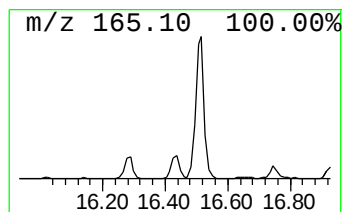
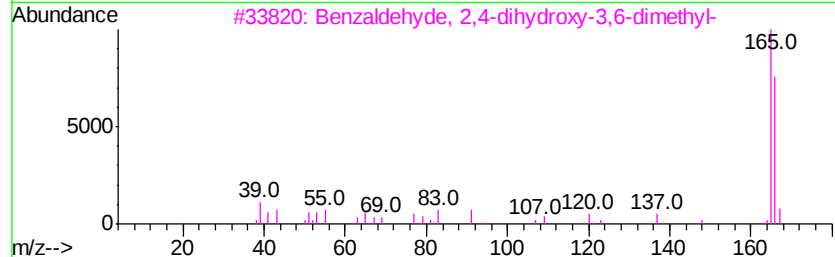
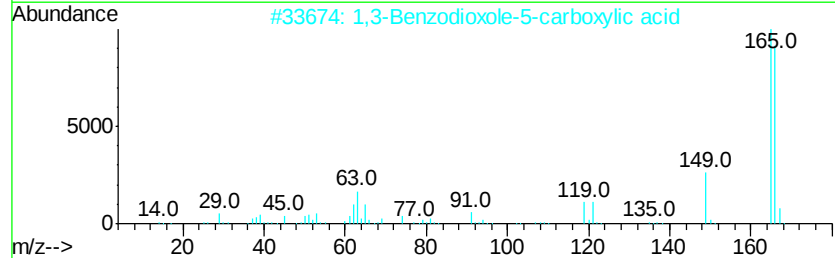
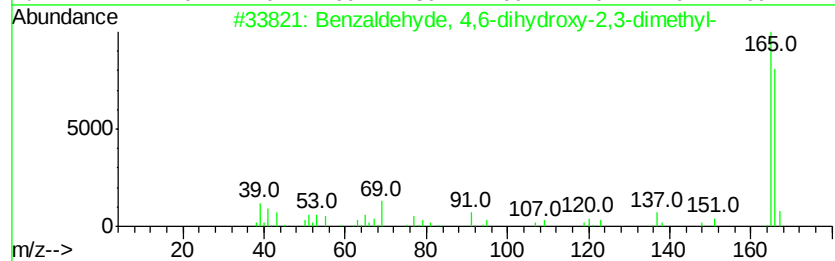
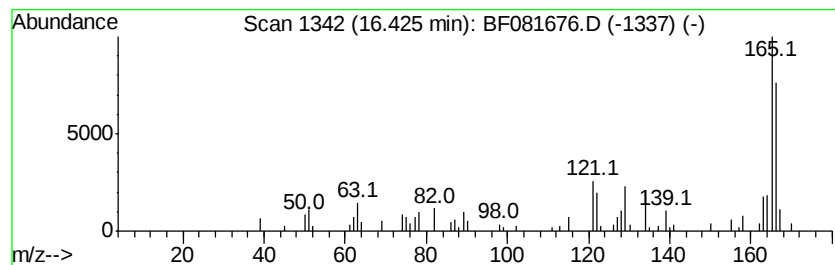
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzaldehyde, 4,6-dihydroxy... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	13.39 ng	256597	Acenaphthene-d10	15.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde, 4,6-dihydroxy-2,3-...	166	C9H10O3	002990-31-0	52
2		1,3-Benzodioxole-5-carboxylic acid	166	C8H6O4	000094-53-1	52
3		Benzaldehyde, 2,4-dihydroxy-3,6-...	166	C9H10O3	034883-14-2	50
4		Fluorene	166	C13H10	000086-73-7	47
5		1H-Phenylene	166	C13H10	000203-80-5	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091715\
 Data File : BF081676.D
 Acq On : 17 Sep 2015 13:17
 Operator : UM/IZ
 Sample : G3669-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-91415-WATER

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.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...	1.22	208.2	ng	2953070	1	8.17	283716	20.0
2-Pentanone, 4-hy...	4.42	22.4	ng	317731	1	8.17	283716	20.0
Benzene, 1-ethyl-...	7.14	14.7	ng	208768	1	8.17	283716	20.0
unknown7.68	7.68	27.0	ng	383305	1	8.17	283716	20.0
Benzene, 1-etheny...	7.83	51.6	ng	731853	1	8.17	283716	20.0
Indane	8.56	27.1	ng	383901	1	8.17	283716	20.0
Indene	8.78	653.5	ng	9270040	1	8.17	283716	20.0
Naphthalene, 1-et...	14.12	9.8	ng	187937	3	15.20	383224	20.0
Naphthalene, 2,6-...	14.27	19.7	ng	377751	3	15.20	383224	20.0
Naphthalene, 2,3-...	14.45	22.5	ng	430315	3	15.20	383224	20.0
Naphthalene, 2,7-...	14.51	10.8	ng	207570	3	15.20	383224	20.0
Naphthalene, 1,5-...	14.71	11.3	ng	216431	3	15.20	383224	20.0
Benzaldehyde, 4,6...	16.43	13.4	ng	256597	3	15.20	383224	20.0