

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043015\
 Data File : BF078868.D
 Acq On : 1 May 2015 1:07
 Operator : TP/IZ
 Sample : G2044-07 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-62S

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-BF043015.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	1.518	27	29	32	rVB	1358320	1452437	36.38%	7.162%
2	1.667	40	42	50	rVB	155049	276262	6.92%	1.362%
3	1.781	50	52	56	rVV2	38517	83123	2.08%	0.410%
4	1.850	56	58	61	rVV	54707	79269	1.99%	0.391%
5	1.907	61	63	97	rVB	2140734	3992671	100.00%	19.688%
6	5.176	347	349	353	rVB	120229	130908	3.28%	0.646%
7	5.667	389	392	394	rBV	251210	257864	6.46%	1.272%
8	6.913	499	501	503	rBV	395135	336620	8.43%	1.660%
9	7.073	513	515	518	rBV	299642	330364	8.27%	1.629%
10	7.370	538	541	543	rBV	530058	465433	11.66%	2.295%
11	7.553	555	557	560	rVB	276154	277917	6.96%	1.370%
12	8.056	599	601	604	rVB	241750	205653	5.15%	1.014%
13	8.948	677	679	681	rBV	761021	656351	16.44%	3.236%
14	10.296	794	797	799	rBV	422151	447829	11.22%	2.208%
15	10.799	838	841	843	rBV	39837	42304	1.06%	0.209%
16	11.131	867	870	872	rBV	542117	711344	17.82%	3.508%
17	12.102	951	955	959	rBV	109420	115416	2.89%	0.569%
18	12.971	1028	1031	1032	rBV	820395	768481	19.25%	3.789%
19	13.005	1032	1034	1036	rVV	361340	450660	11.29%	2.222%
20	13.062	1037	1039	1042	rVB	118666	113475	2.84%	0.560%
21	13.394	1066	1068	1070	rVB	28596	39982	1.00%	0.197%
22	13.599	1083	1086	1088	rBV	77958	93659	2.35%	0.462%
23	13.634	1088	1089	1091	rVV	107031	93923	2.35%	0.463%
24	13.691	1091	1094	1096	rVV	43024	51861	1.30%	0.256%
25	13.737	1096	1098	1099	rVV	118285	142279	3.56%	0.702%
26	13.759	1099	1100	1104	rVB	48625	71403	1.79%	0.352%
27	13.897	1110	1112	1117	rBV	31182	49151	1.23%	0.242%
28	13.977	1117	1119	1123	rBV2	59867	102955	2.58%	0.508%
29	14.285	1144	1146	1148	rBV	51891	67511	1.69%	0.333%
30	14.331	1148	1150	1153	rVV3	33370	54636	1.37%	0.269%
31	14.388	1153	1155	1158	rVB2	56456	77798	1.95%	0.384%
32	14.480	1160	1163	1166	rBV	859346	833140	20.87%	4.108%
33	14.560	1168	1170	1172	rBV2	29597	43843	1.10%	0.216%
34	14.662	1176	1179	1184	rBV2	26879	74372	1.86%	0.367%

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35	14.765	1184	1188	1190	rBV	651062	889246	22.27%	4.385%
36	14.822	1190	1193	1195	rVV	15334	43456	1.09%	0.214%
37	14.914	1199	1201	1203	rBV	43232	60942	1.53%	0.301%
38	14.971	1203	1206	1208	rBV	402215	586592	14.69%	2.892%
39	15.040	1208	1212	1214	rBV3	40328	65913	1.65%	0.325%
40	15.154	1219	1222	1223	rBV3	32487	66157	1.66%	0.326%
41	15.177	1223	1224	1226	rVB	83081	77277	1.94%	0.381%
42	15.268	1230	1232	1233	rBV	39515	56033	1.40%	0.276%
43	15.302	1233	1235	1237	rBV	100849	114276	2.86%	0.563%
44	15.417	1243	1245	1247	rBV	49643	57940	1.45%	0.286%
45	15.462	1247	1249	1251	rVV	38827	54856	1.37%	0.270%
46	15.805	1277	1279	1282	rBV	63266	80160	2.01%	0.395%
47	15.943	1289	1291	1294	rBV2	54348	96208	2.41%	0.474%
48	15.988	1294	1295	1298	rVV2	61793	122961	3.08%	0.606%
49	16.068	1301	1302	1305	rVB	45773	56620	1.42%	0.279%
50	16.263	1316	1319	1321	rBV2	928170	1373192	34.39%	6.771%
51	16.297	1321	1322	1324	rVV	330903	255400	6.40%	1.259%
52	16.388	1328	1330	1332	rBV2	50898	66255	1.66%	0.327%
53	16.754	1360	1362	1364	rVB	78916	88087	2.21%	0.434%
54	16.800	1364	1366	1368	rBV	41033	43807	1.10%	0.216%
55	17.531	1427	1430	1436	rBV	351955	833151	20.87%	4.108%
56	17.646	1438	1440	1442	rVB	73989	101180	2.53%	0.499%
57	17.851	1455	1458	1461	rVB	222579	294990	7.39%	1.455%
58	17.908	1461	1463	1466	rBV	262650	397053	9.94%	1.958%
59	17.988	1467	1470	1472	rBV	525195	838402	21.00%	4.134%
60	19.486	1598	1601	1606	rBV2	135861	281238	7.04%	1.387%
61	19.931	1637	1640	1647	rVB	129500	287704	7.21%	1.419%

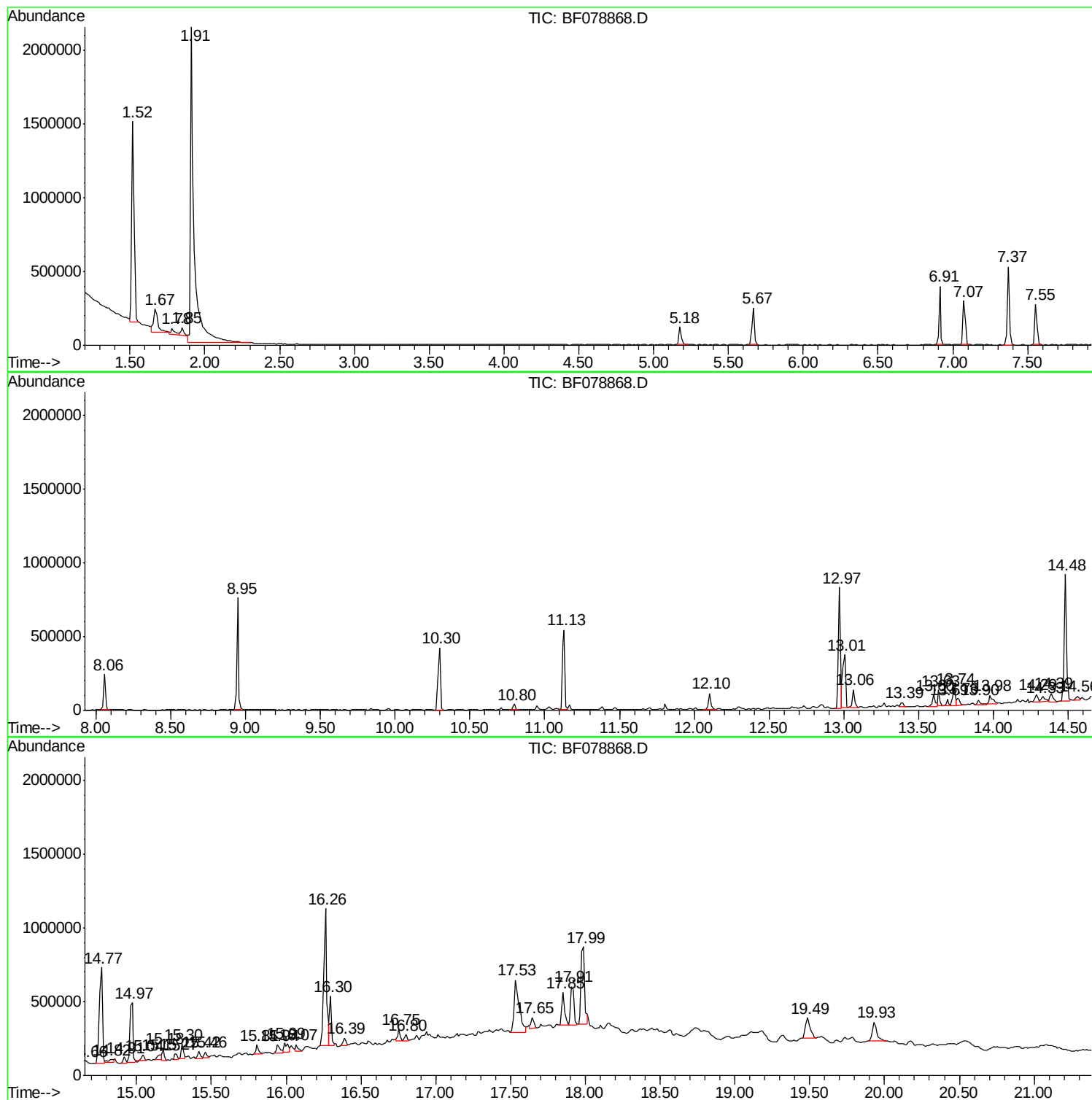
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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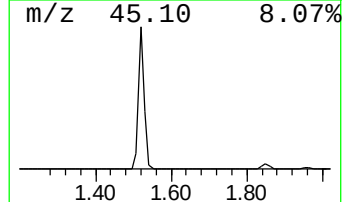
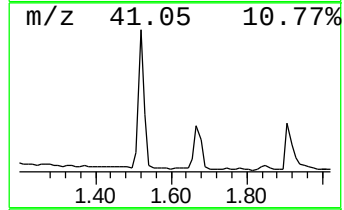
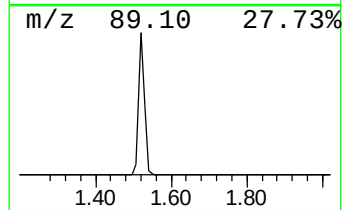
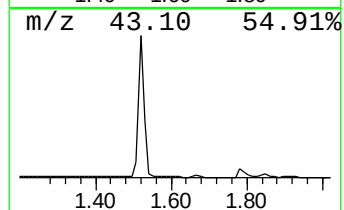
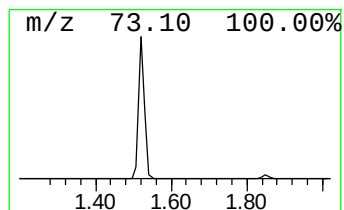
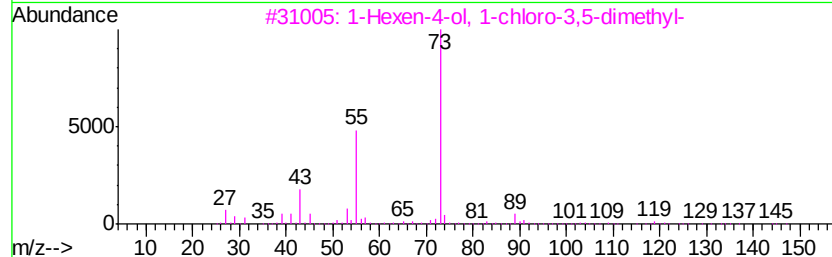
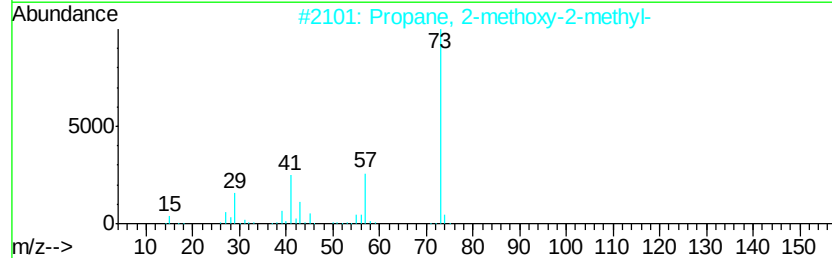
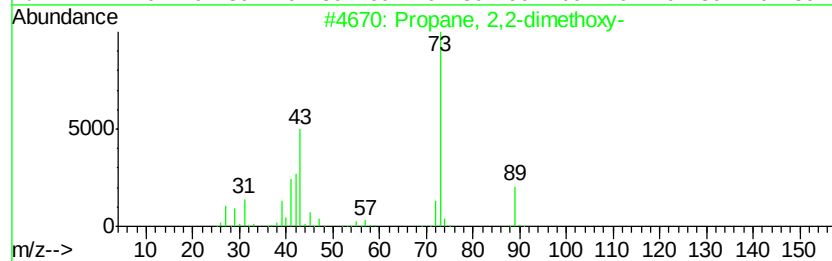
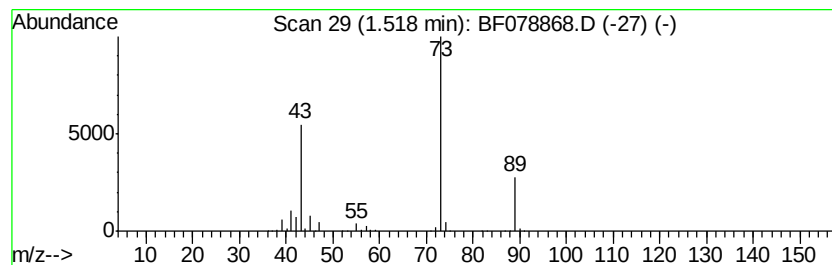
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	62.41 ng	1452440	1,4-Dichlorobenzene-d4	7.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
4		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
5		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9



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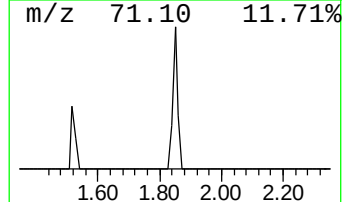
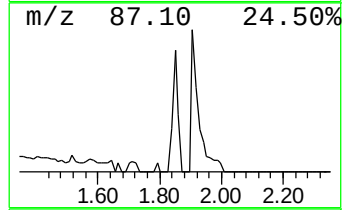
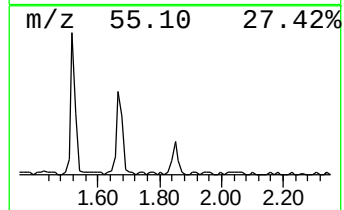
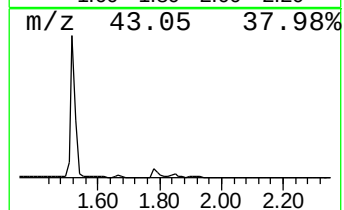
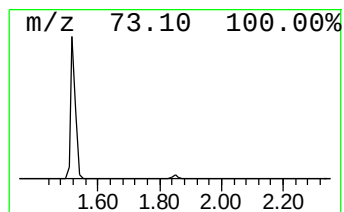
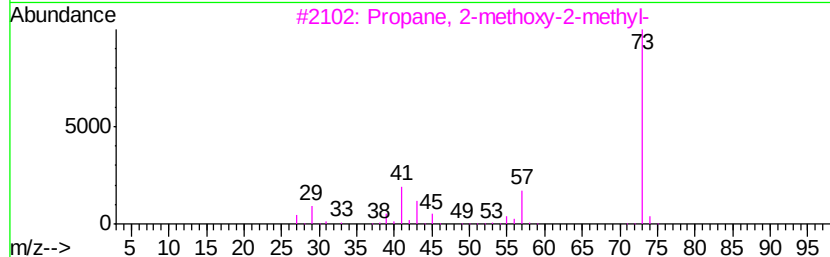
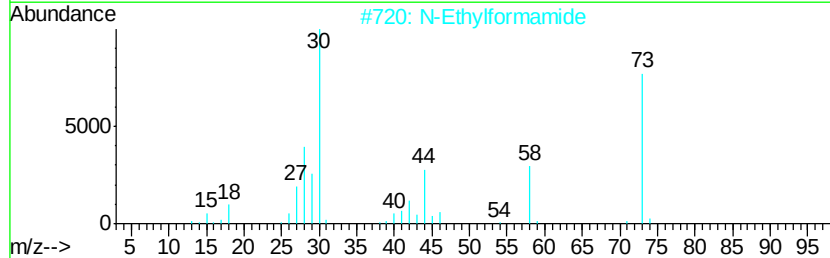
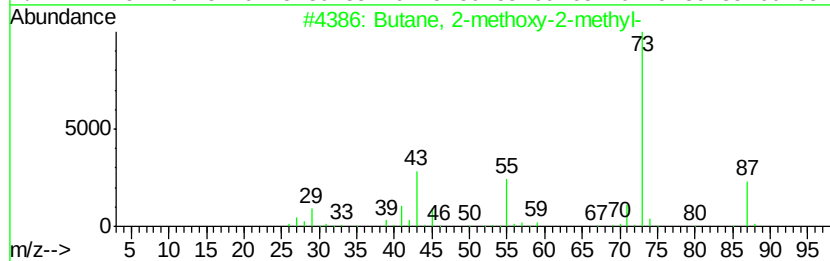
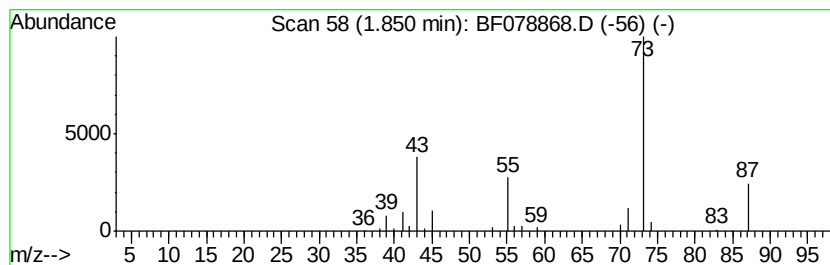
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.85	3.41 ng	79269	1,4-Dichlorobenzene-d4	7.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		N-Ethylformamide	73	C3H7NO	000627-45-2	9
3		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	5



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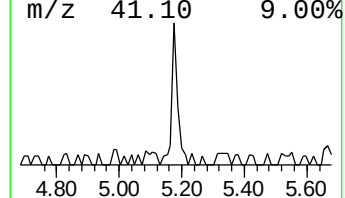
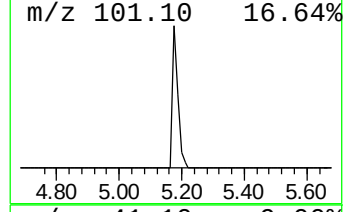
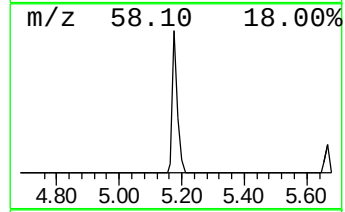
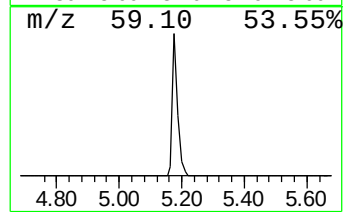
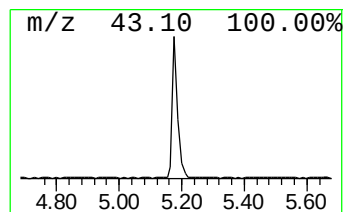
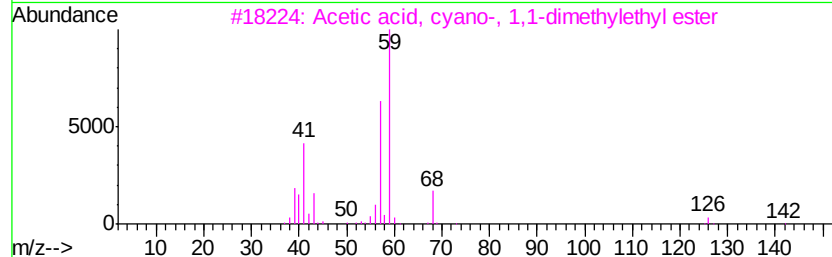
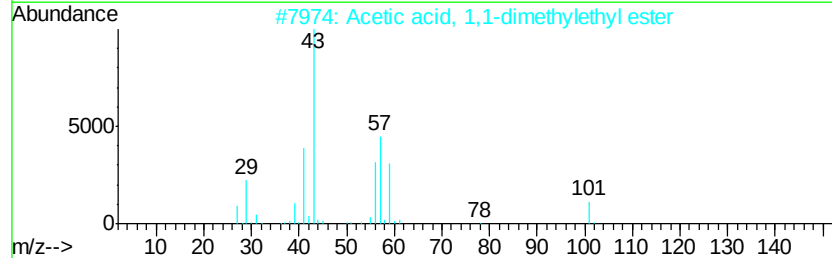
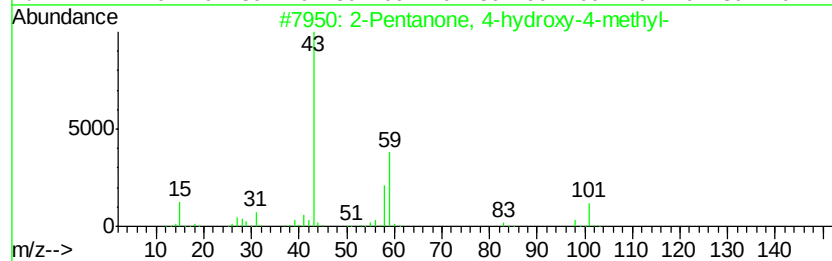
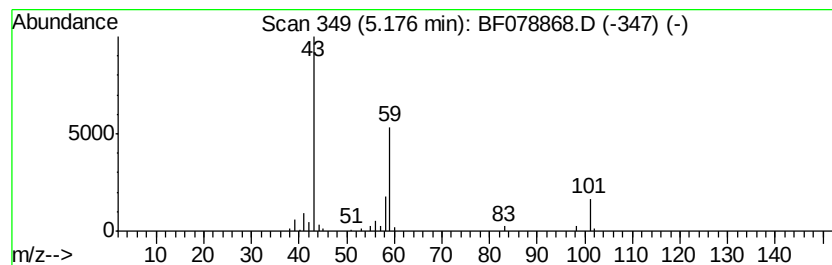
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	5.63 ng	130908	1,4-Dichlorobenzene-d4	7.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	10



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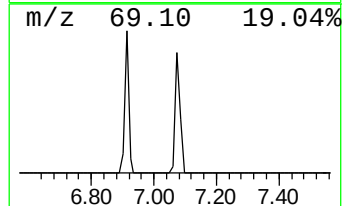
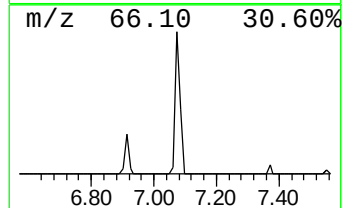
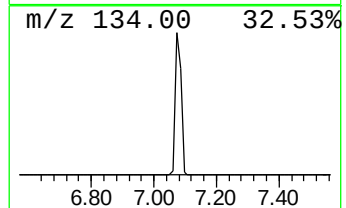
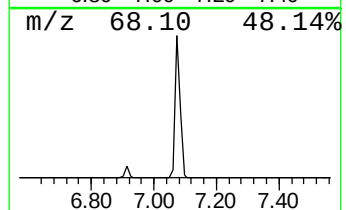
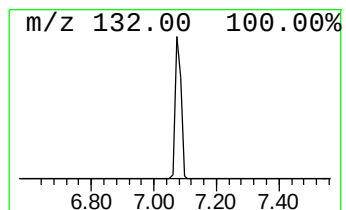
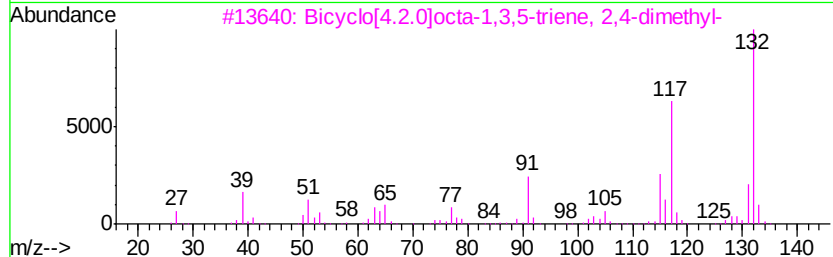
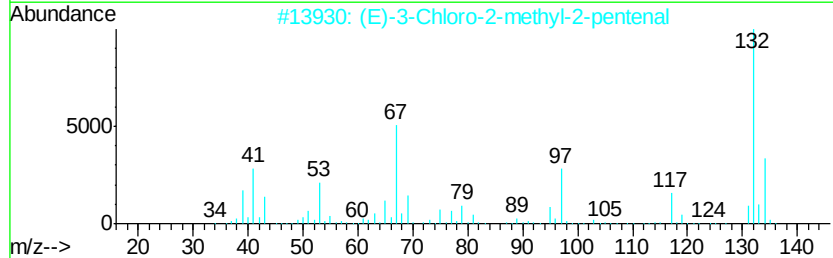
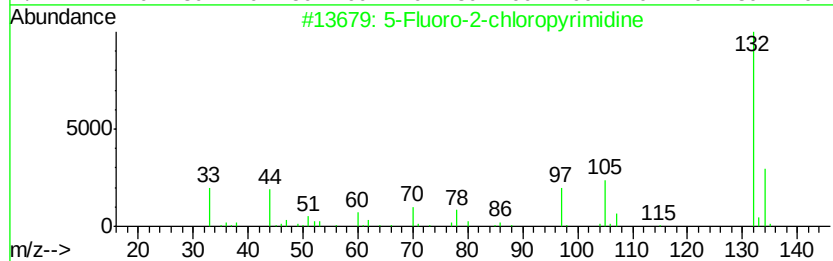
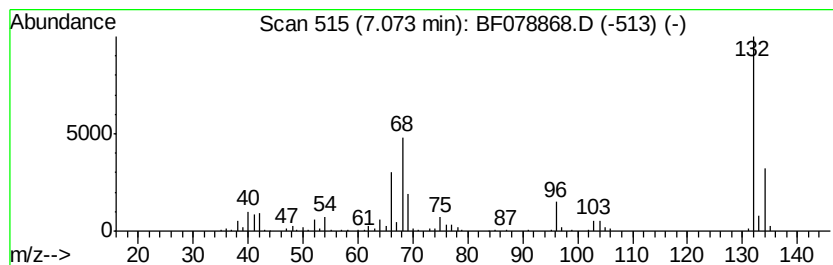
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown7.07 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	14.20 ng	330364	1,4-Dichlorobenzene-d4	7.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
3		Bicyclo[4.2.0]octa-1,3,5-triene, ...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9



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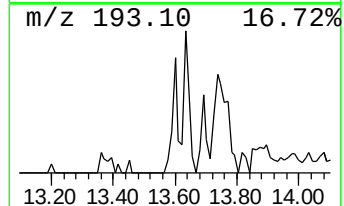
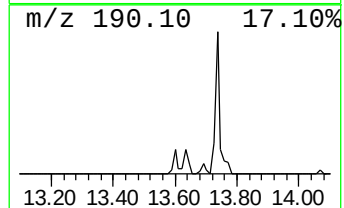
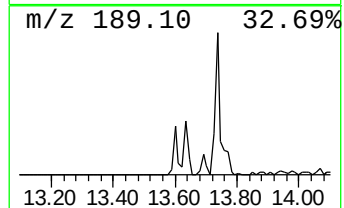
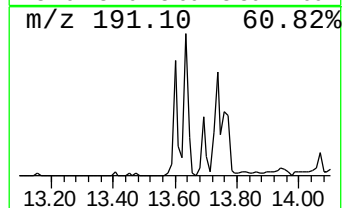
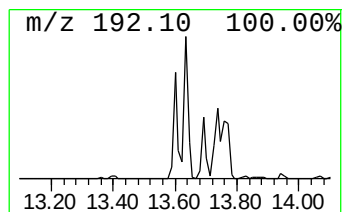
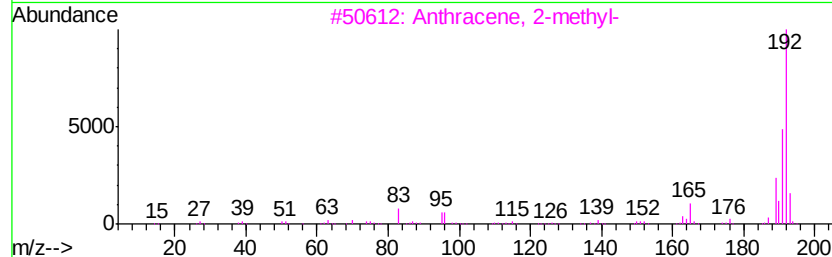
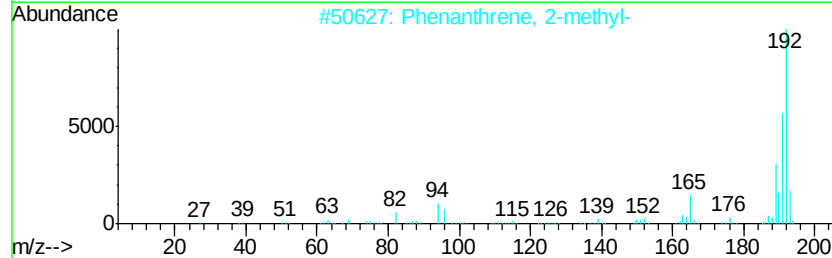
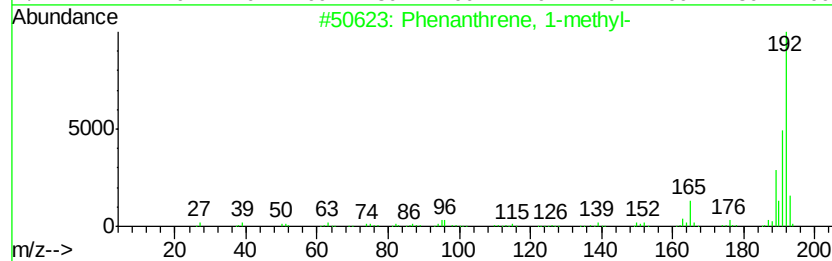
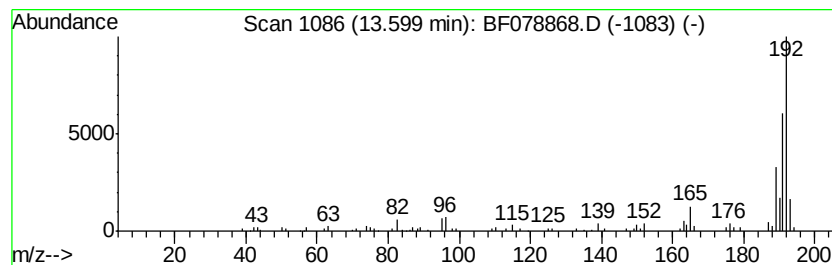
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.60	2.44 ng	93659	Phenanthrene-d10	12.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
5	1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF043015\
Data File : BF078868.D
Acq On : 1 May 2015 1:07
Operator : TP/IZ
Sample : G2044-07 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-62S

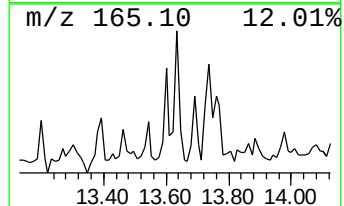
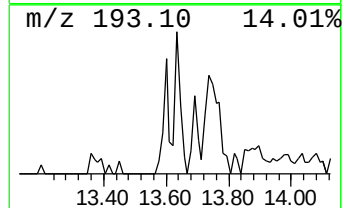
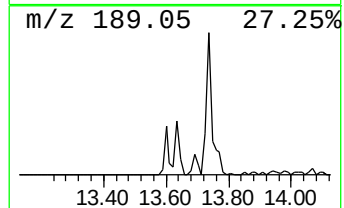
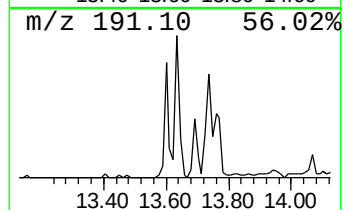
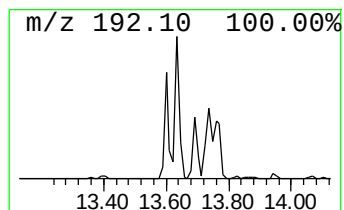
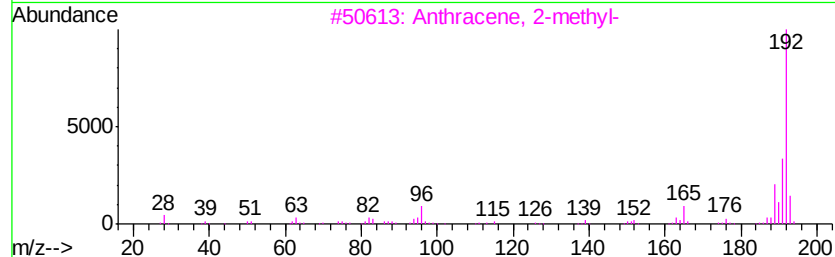
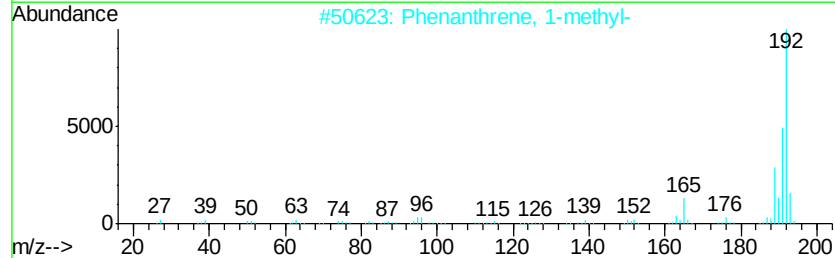
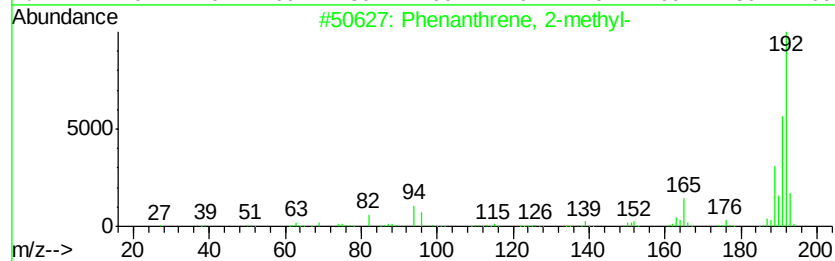
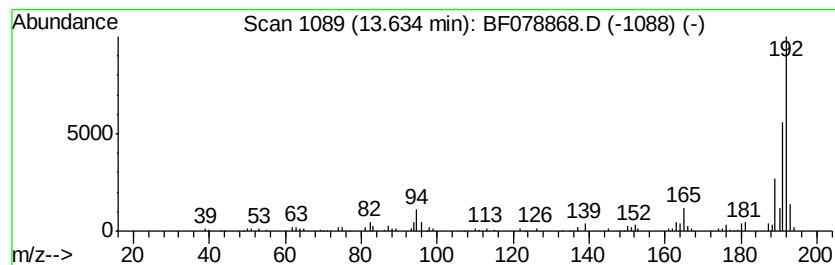
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	2.44 ng	93923	Phenanthrene-d10	12.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
5		1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043015\
Data File : BF078868.D
Acq On : 1 May 2015 1:07
Operator : TP/IZ
Sample : G2044-07 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-62S

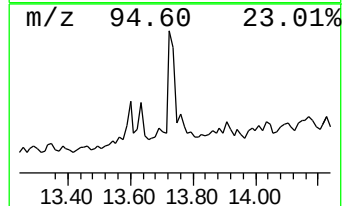
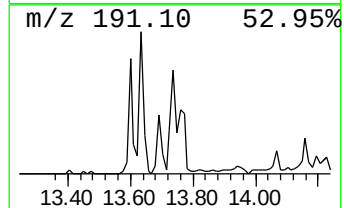
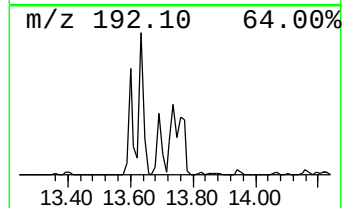
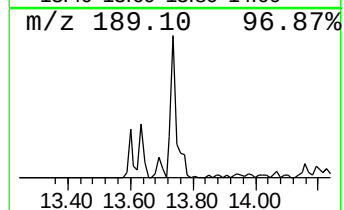
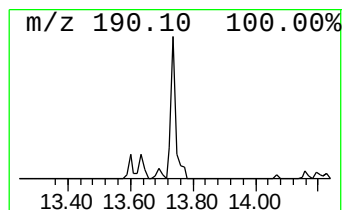
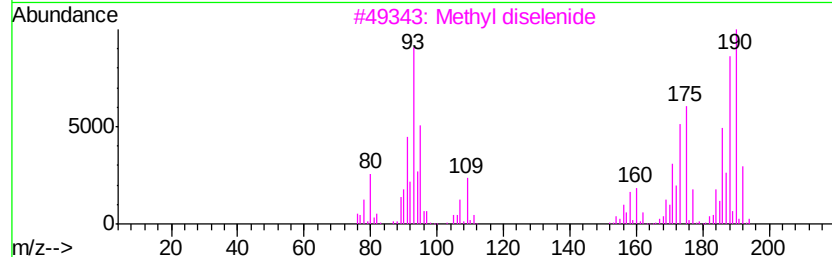
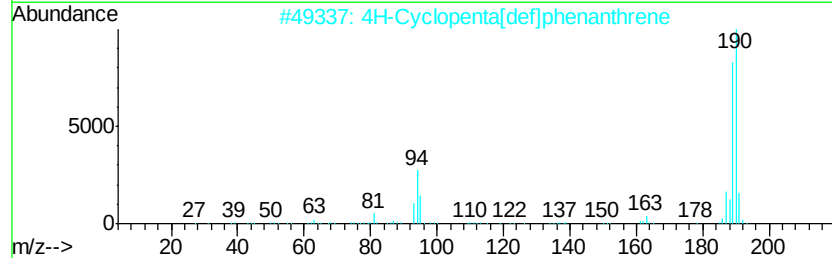
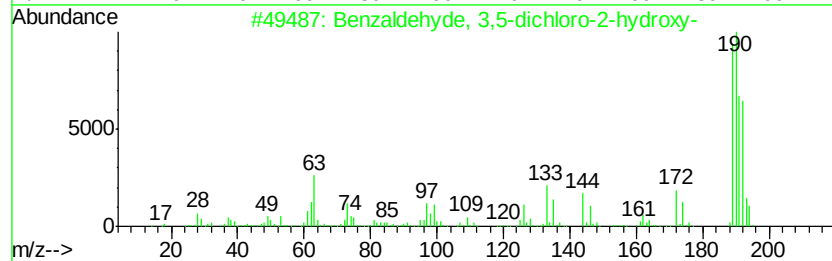
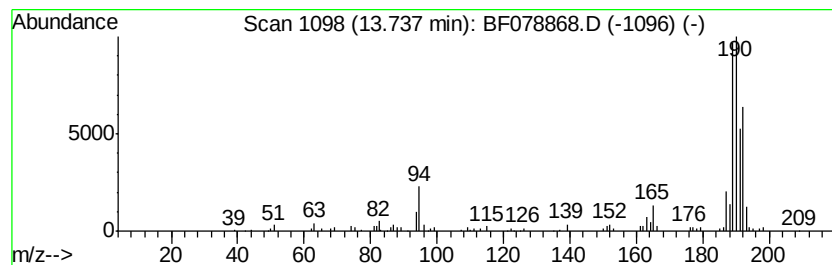
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzaldehyde, 3,5-dichloro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.74	3.70 ng	142279	Phenanthrene-d10	12.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
3			Methyl diselenide	190	C2H6Se2	007101-31-7	41
4			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	38
5			8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043015\
Data File : BF078868.D
Acq On : 1 May 2015 1:07
Operator : TP/IZ
Sample : G2044-07 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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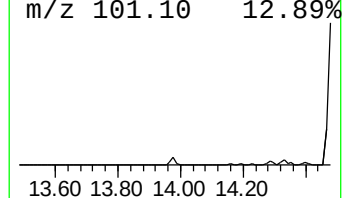
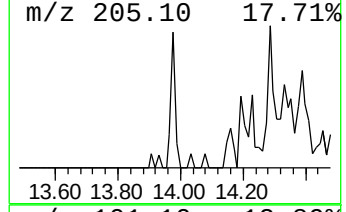
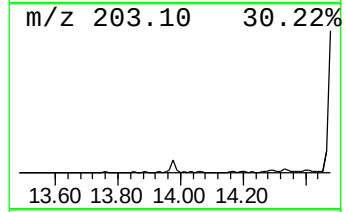
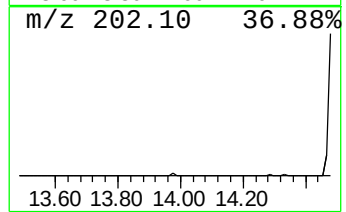
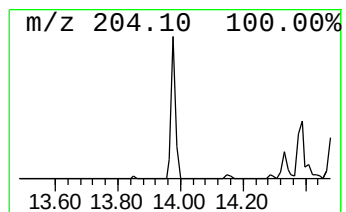
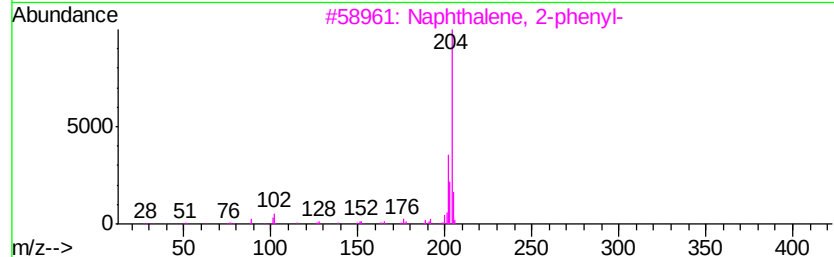
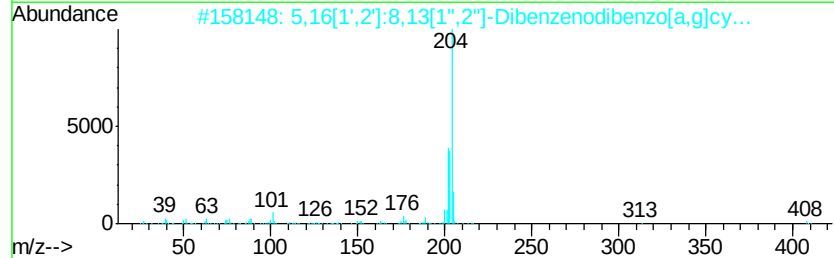
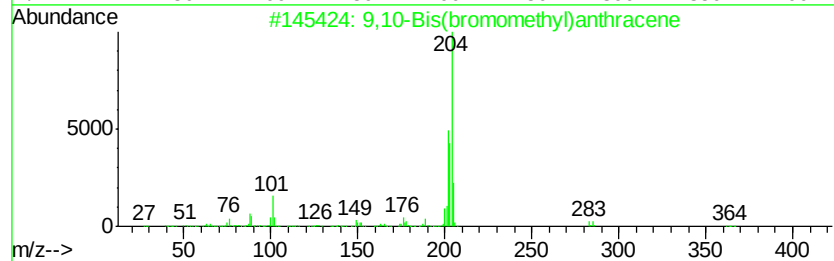
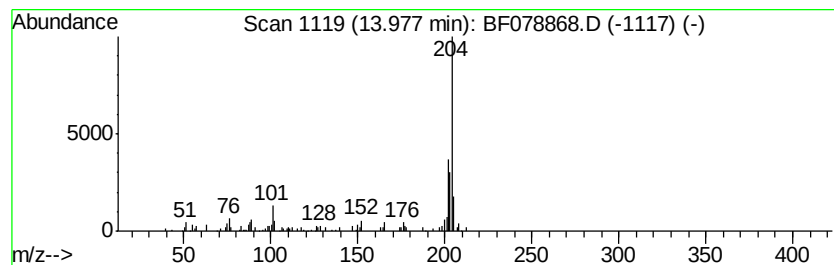
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9,10-Bis(bromomethyl)anthra... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	2.68 ng	102955	Phenanthrene-d10	12.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	90
2		5,16[1',2'] : 8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	81
4		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81
5		2-Phenyl naphthalene	204	C16H12	035465-71-5	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043015\
Data File : BF078868.D
Acq On : 1 May 2015 1:07
Operator : TP/IZ
Sample : G2044-07 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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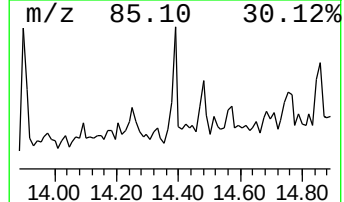
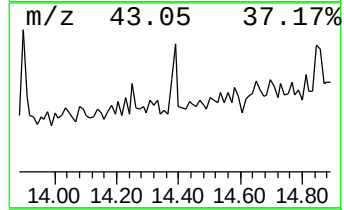
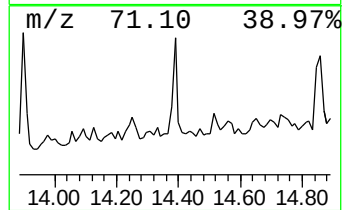
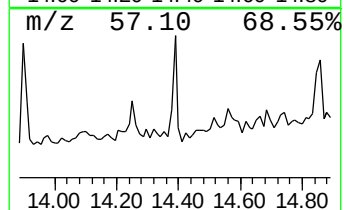
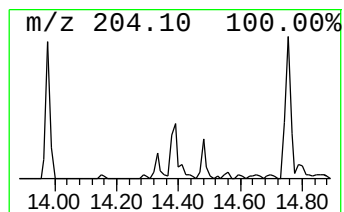
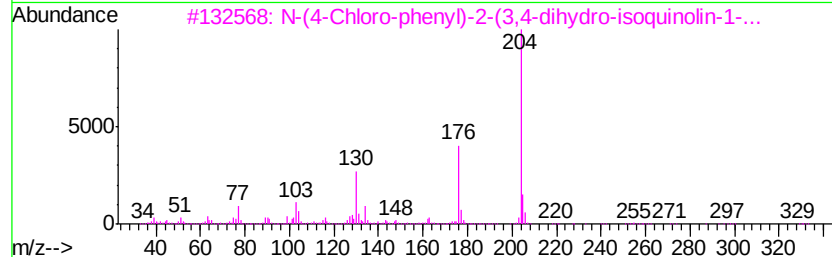
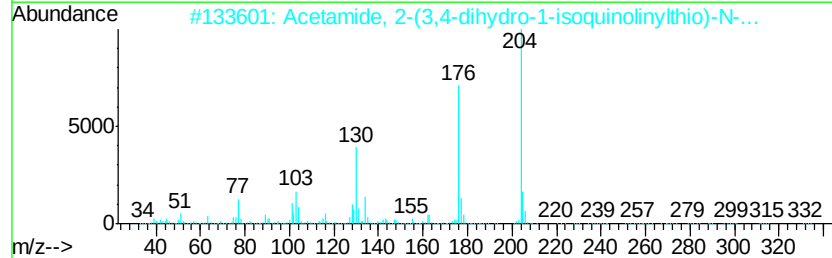
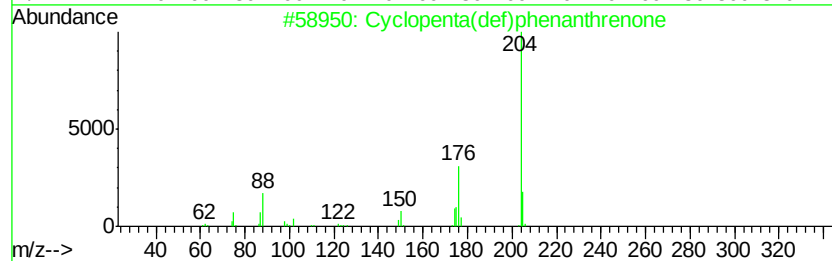
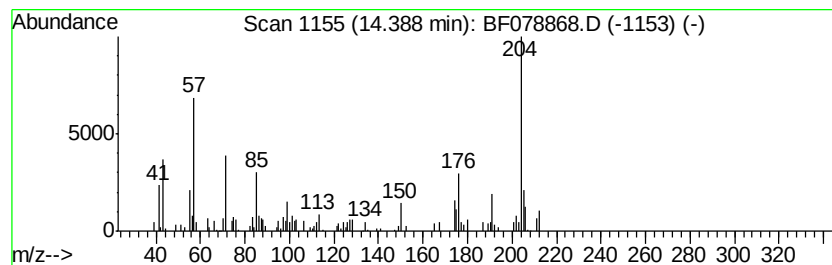
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopenta(def)phenanthrenone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	2.02 ng	77798	Phenanthrene-d10	12.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	60
2		Acetamide, 2-(3,4-dihydro-1-isoq...	332	C17H14F2N2OS	1000270-76-1	43
3		N-(4-Chloro-phenyl)-2-(3,4-dihvd...	330	C17H15ClN2OS	1000296-34-3	43
4		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	38
5		5H-Dibenzo[a,d]cycloheptene, 5-m...	204	C16H12	002975-79-3	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043015\
Data File : BF078868.D
Acq On : 1 May 2015 1:07
Operator : TP/IZ
Sample : G2044-07 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

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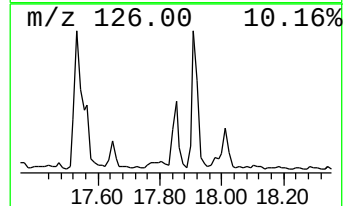
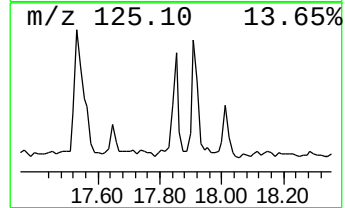
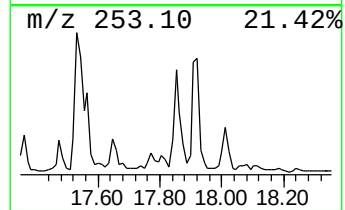
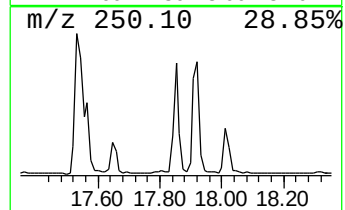
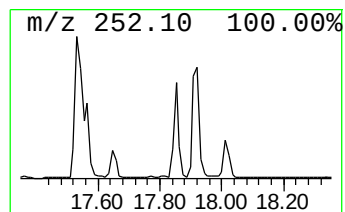
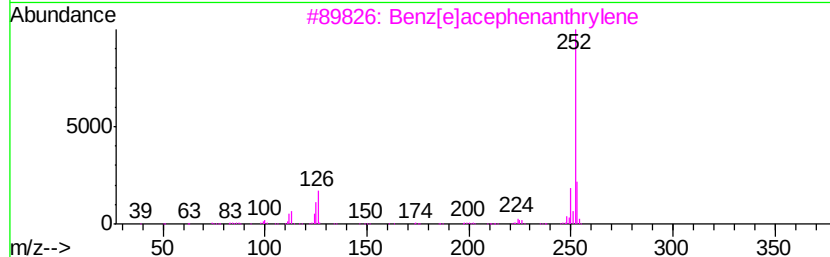
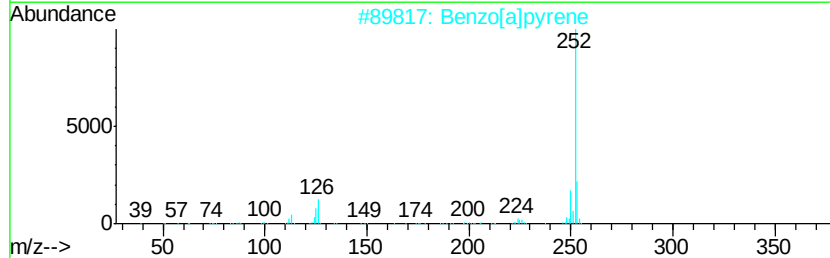
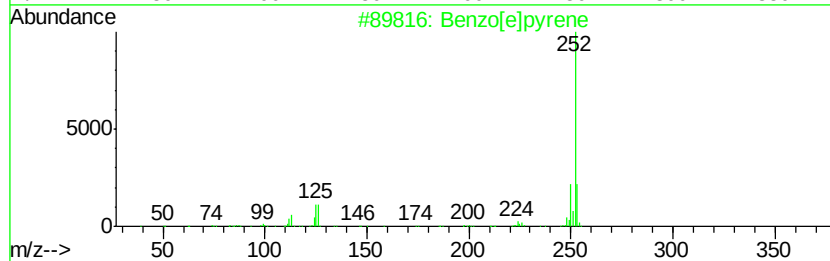
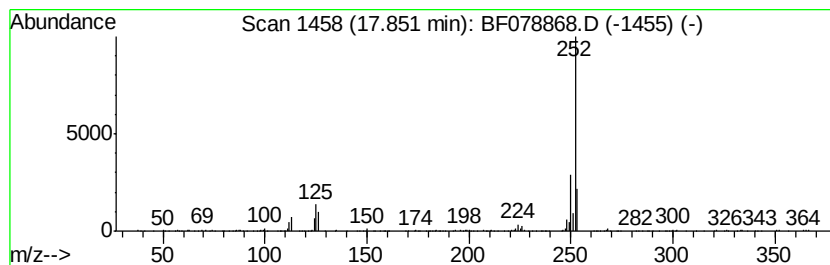
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.85	7.04 ng	294990	Perylene-d12	17.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[a]pyrene	252	C20H12	000050-32-8	96
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	95
4		Benzo[i]lfluoranthene	252	C20H12	000205-82-3	95
5		Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF043015\
Data File : BF078868.D
Acq On : 1 May 2015 1:07
Operator : TP/IZ
Sample : G2044-07 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
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Butane, 2-methoxy...	1.85	3.4	ng	79269	1	7.37	465433	20.0
2-Pentanone, 4-hy...	5.18	5.6	ng	130908	1	7.37	465433	20.0
unknown7.07	7.07	14.2	ng	330364	1	7.37	465433	20.0
Phenanthrene, 1-m...	13.60	2.4	ng	93659	4	12.97	768481	20.0
Phenanthrene, 2-m...	13.63	2.4	ng	93923	4	12.97	768481	20.0
Benzaldehyde, 3,5...	13.74	3.7	ng	142279	4	12.97	768481	20.0
9,10-Bis(bromomet...	13.98	2.7	ng	102955	4	12.97	768481	20.0
Cyclopenta(def)ph...	14.39	2.0	ng	77798	4	12.97	768481	20.0
Benzo[e]pyrene	17.85	7.0	ng	294990	6	17.99	838402	20.0