

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
 Data File : BF078974.D
 Acq On : 6 May 2015 21:47
 Operator : TP/IZ
 Sample : G2035-17
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-15-SS

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	1.495	24	27	30	rBV	5680480	6937928	100.00%	21.094%
2	1.633	36	39	49	rVB	149399	271286	3.91%	0.825%
3	1.770	49	51	53	rVV	117368	159852	2.30%	0.486%
4	1.815	53	55	60	rVV	172544	248582	3.58%	0.756%
5	1.896	60	62	83	rVB	767409	1227299	17.69%	3.731%
6	5.153	345	347	356	rVB	309891	396305	5.71%	1.205%
7	5.644	388	390	393	rBV	1532211	1956159	28.20%	5.947%
8	6.902	498	500	503	rBV	2054443	2009970	28.97%	6.111%
9	7.062	511	514	516	rBV	2368261	2181750	31.45%	6.633%
10	7.347	537	539	541	rBV	444958	415138	5.98%	1.262%
11	7.542	553	556	558	rVB	1247348	1577571	22.74%	4.796%
12	8.045	597	600	602	rBV	1084082	1345694	19.40%	4.091%
13	8.925	675	677	680	rBV	702222	611003	8.81%	1.858%
14	10.273	792	795	797	rBV	2610794	2303046	33.20%	7.002%
15	10.776	837	839	841	rBV	107025	96360	1.39%	0.293%
16	11.108	865	868	870	rVB	624000	696747	10.04%	2.118%
17	12.034	946	949	951	rBV	182193	195740	2.82%	0.595%
18	12.091	951	954	956	rBV2	1316908	1763770	25.42%	5.362%
19	12.948	1027	1029	1031	rBV	677591	758367	10.93%	2.306%
20	14.365	1151	1153	1157	rVV	77499	76309	1.10%	0.232%
21	14.457	1159	1161	1166	rVB	54610	87919	1.27%	0.267%
22	14.720	1182	1184	1187	rBV2	64465	167735	2.42%	0.510%
23	14.765	1187	1188	1190	rVV	290104	267845	3.86%	0.814%
24	14.834	1190	1194	1195	rVV2	52606	122102	1.76%	0.371%
25	14.868	1195	1197	1198	rVV	58151	81911	1.18%	0.249%
26	14.948	1202	1204	1207	rVB	2468520	2541727	36.64%	7.728%
27	15.657	1263	1266	1268	rBV	166890	214466	3.09%	0.652%
28	16.114	1304	1306	1309	rBV	305517	395453	5.70%	1.202%
29	16.251	1315	1318	1320	rBV	901395	1013932	14.61%	3.083%
30	16.537	1341	1343	1346	rVV	95621	134011	1.93%	0.407%
31	16.948	1377	1379	1382	rBV	111856	166621	2.40%	0.507%
32	17.234	1402	1404	1408	rBV2	113654	225451	3.25%	0.685%
33	17.348	1411	1414	1419	rVB	95632	242310	3.49%	0.737%
34	17.737	1446	1448	1452	rBV2	89851	218344	3.15%	0.664%

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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-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	18.023	1470	1473	1481	rVB	805765	1265002	18.23%	3.846%
36	18.149	1482	1484	1486	rBV	64852	103179	1.49%	0.314%
37	18.606	1522	1524	1526	rBV	75667	111343	1.60%	0.339%
38	20.172	1658	1661	1668	rVB4	106790	302688	4.36%	0.920%

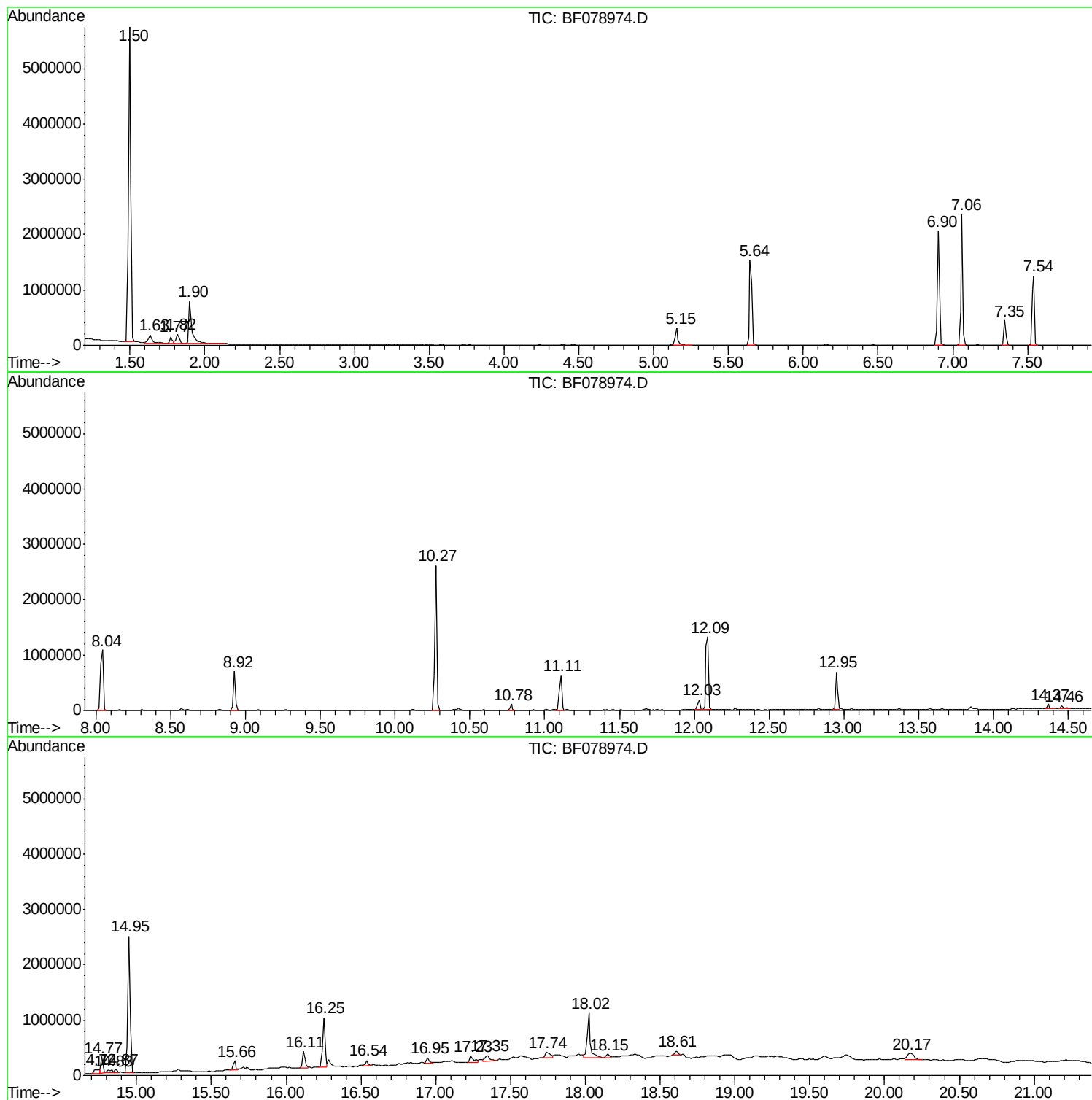
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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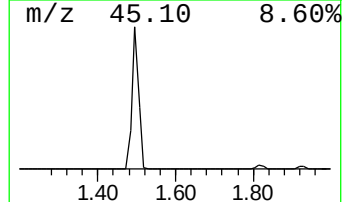
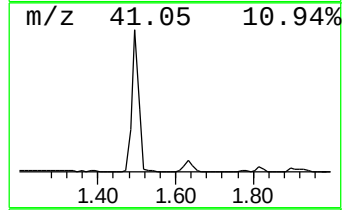
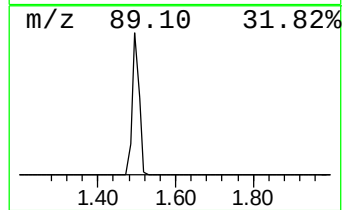
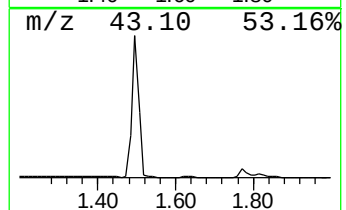
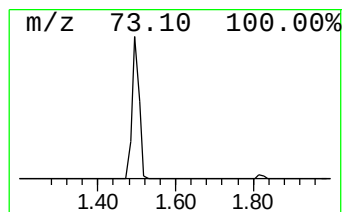
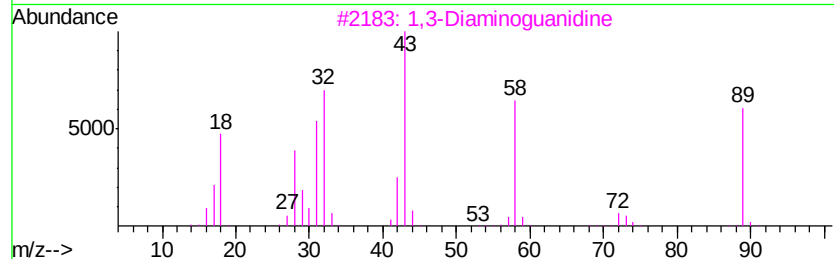
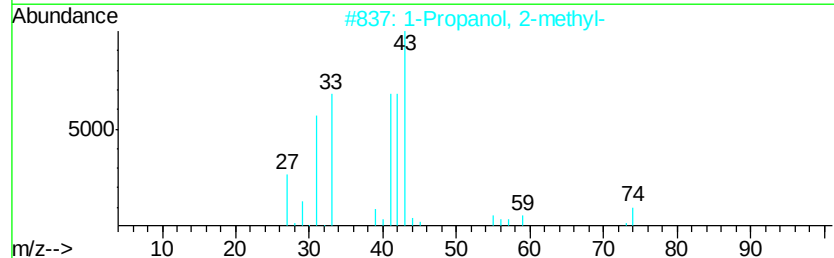
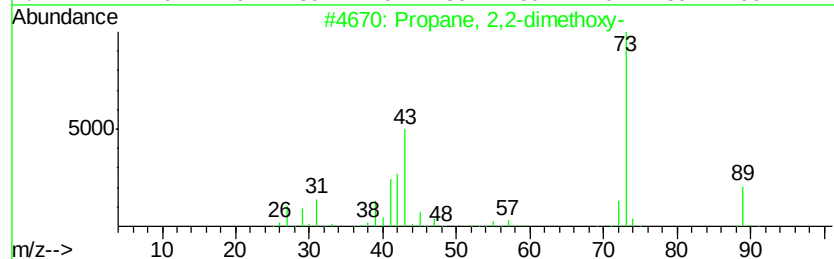
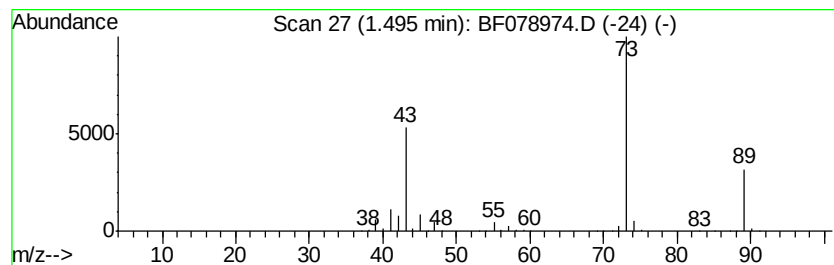
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.50	334.25 ng	6937930	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
4		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
5		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9



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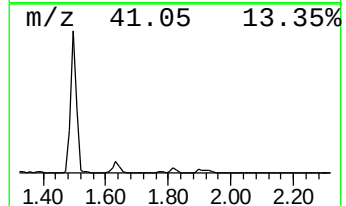
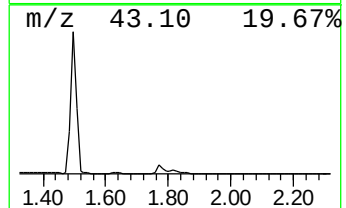
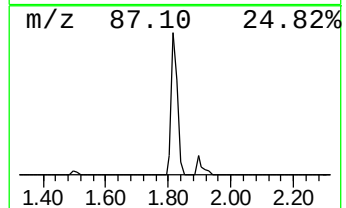
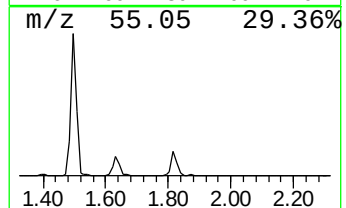
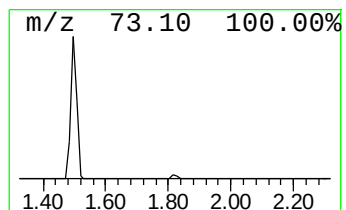
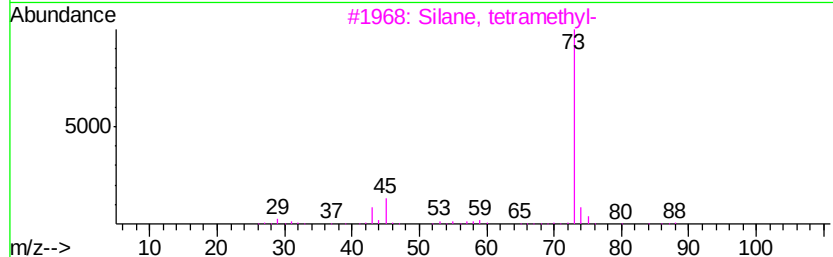
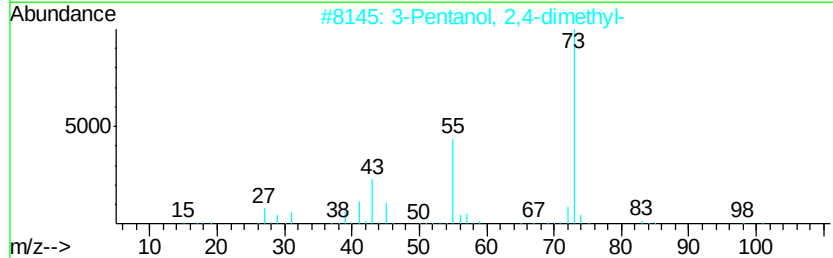
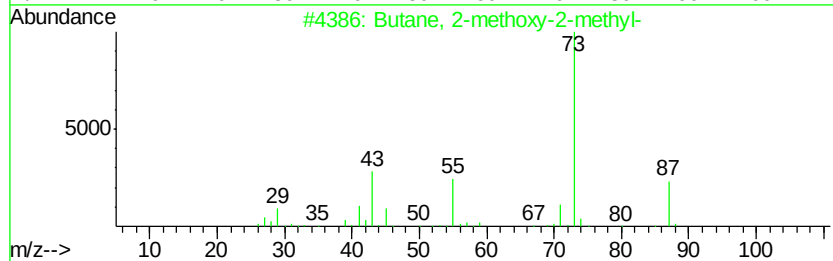
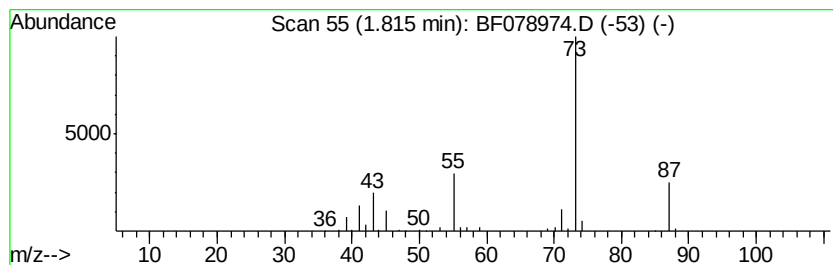
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.82	11.98 ng	248582	1,4-Dichlorobenzene-d4	7.35

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	23
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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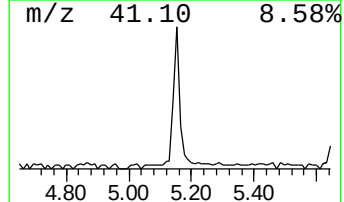
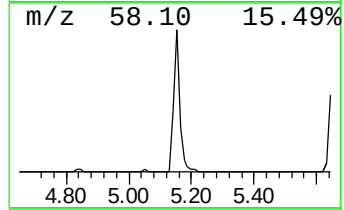
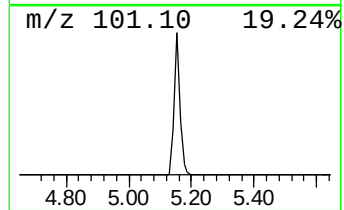
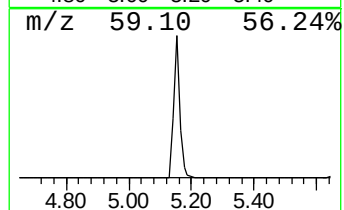
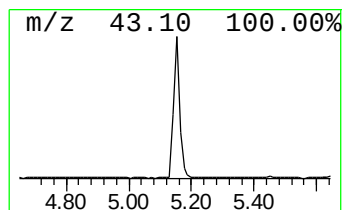
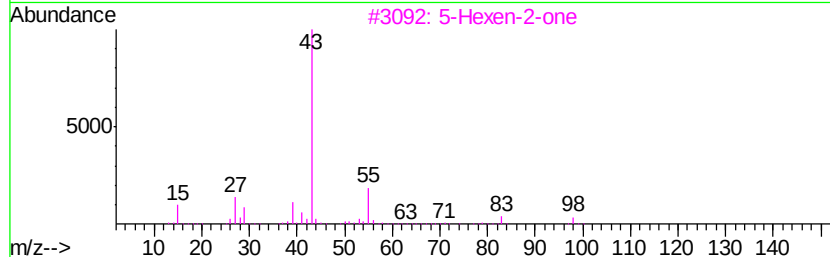
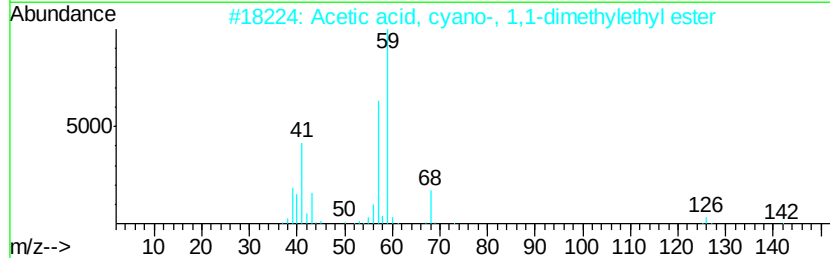
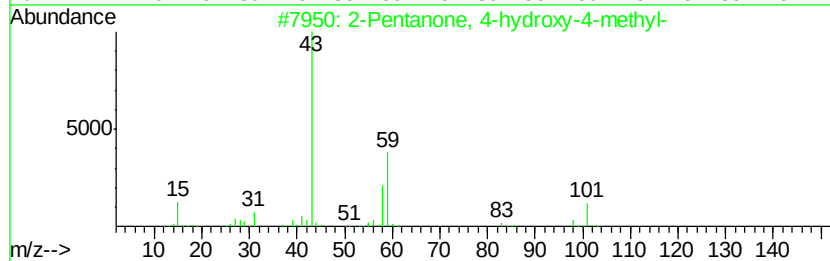
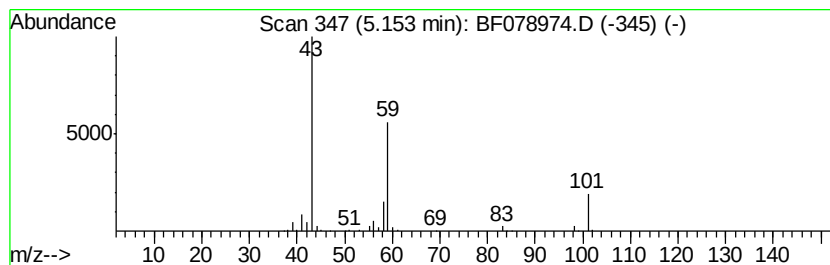
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	19.09 ng	396305	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		t-Butyl ethyl ether	102	C6H14O	1000118-18-4	9



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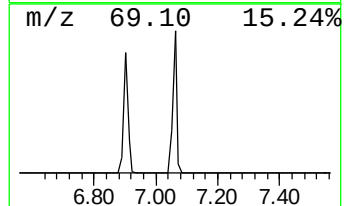
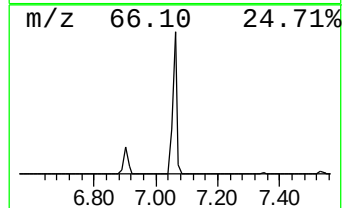
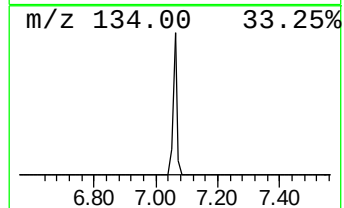
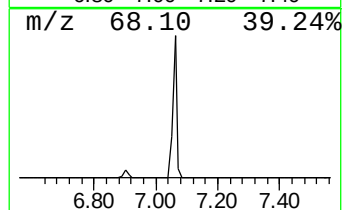
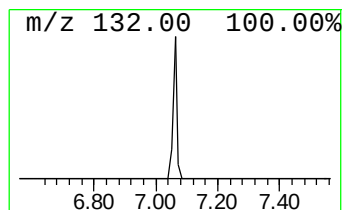
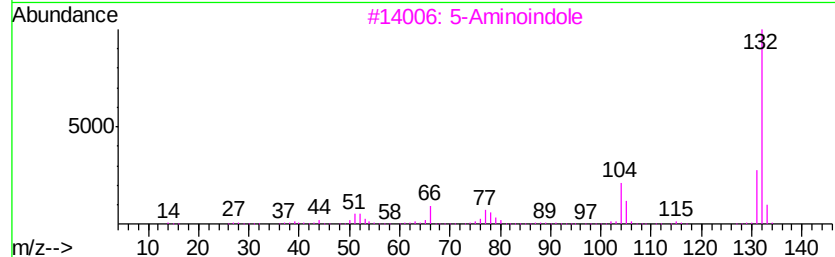
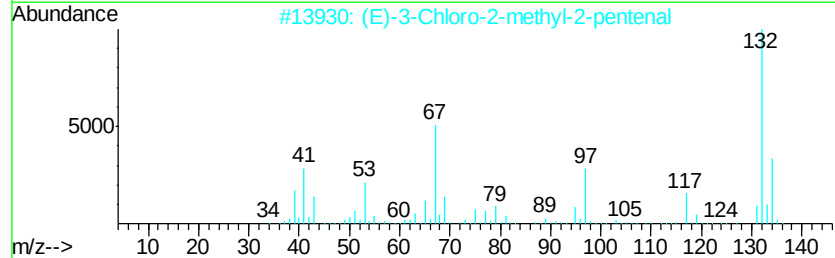
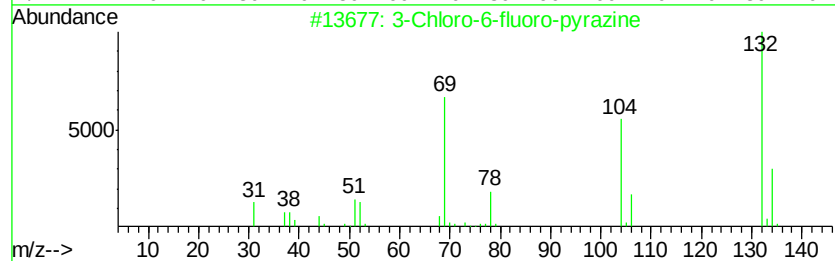
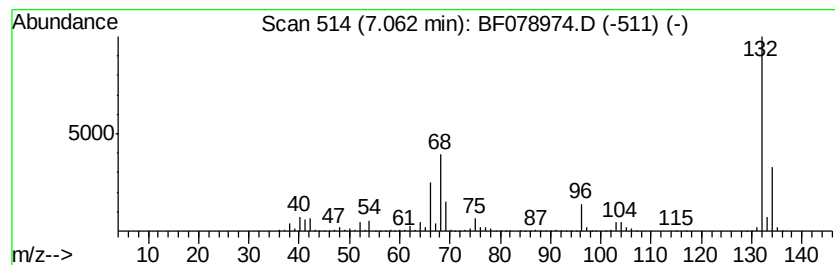
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown7.06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	105.11 ng	2181750	1,4-Dichlorobenzene-d4	7.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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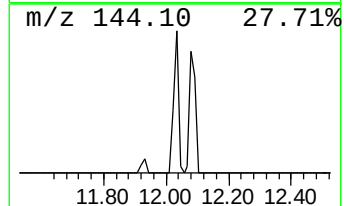
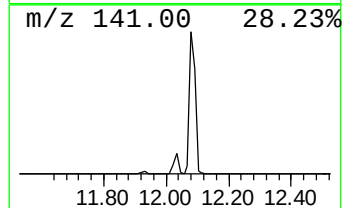
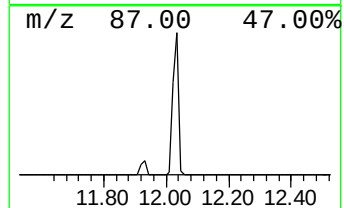
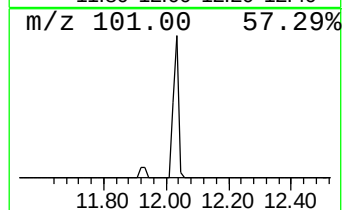
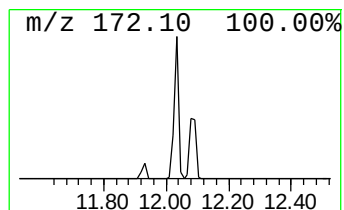
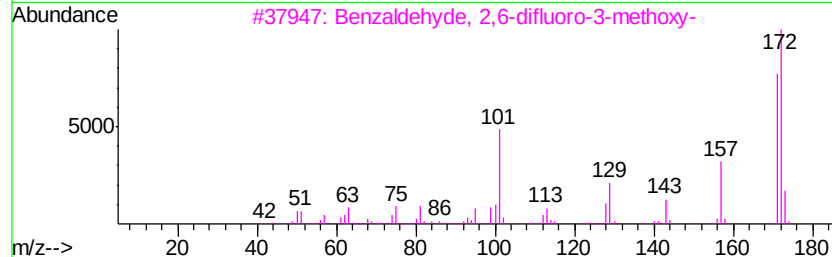
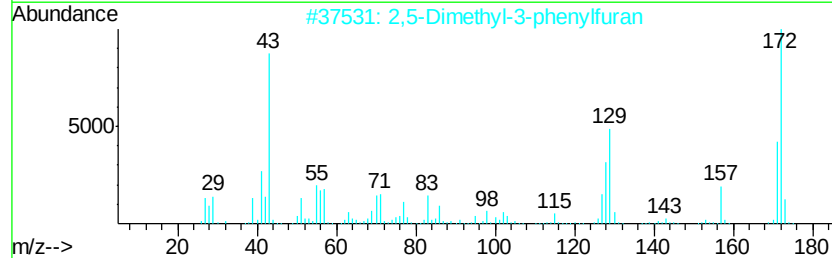
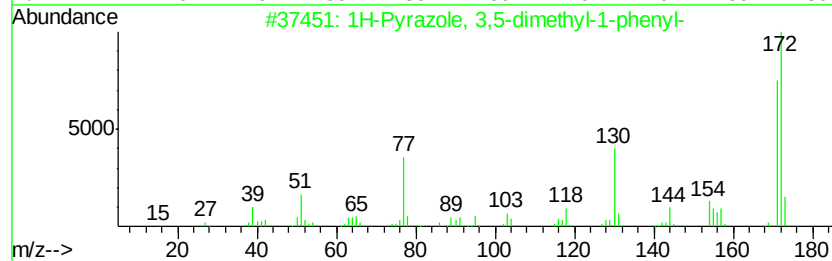
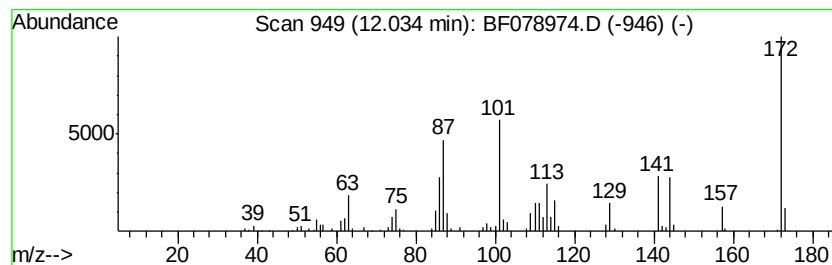
Instrument :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown12.03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.03	5.16 ng	195740	Phenanthrene-d10	12.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrazole, 3,5-dimethyl-1-phenyl-	172	C11H12N2	001131-16-4	27
2	2,5-Dimethyl-3-phenylfuran	172	C12H12O	019842-57-0	27
3	Benzaldehyde, 2,6-difluoro-3-met...	172	C8H6F2O2	149949-30-4	23
4	Menadione	172	C11H8O2	000058-27-5	22
5	peri-Naphtho-1,3-dioxin	172	C11H8O2	000204-11-5	16



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
Data File : BF078974.D
Acq On : 6 May 2015 21:47
Operator : TP/IZ
Sample : G2035-17
Misc :
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Instrument :
BNA_F
ClientSampleId :
TP-15-SS

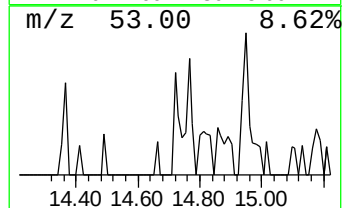
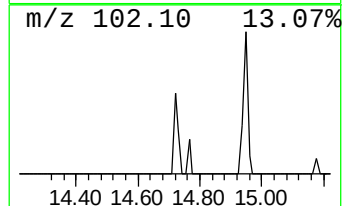
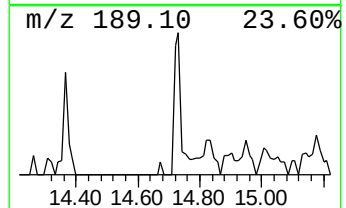
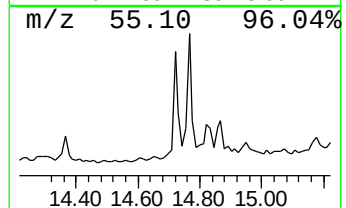
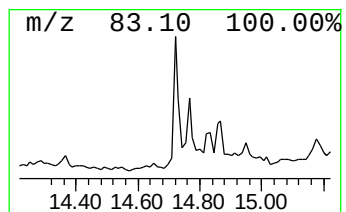
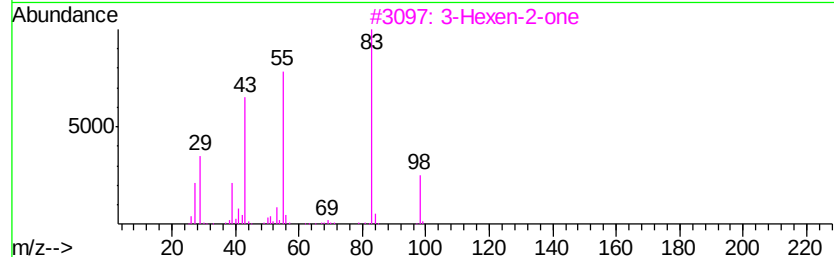
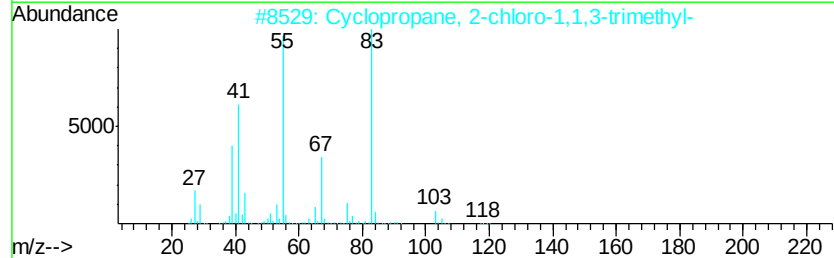
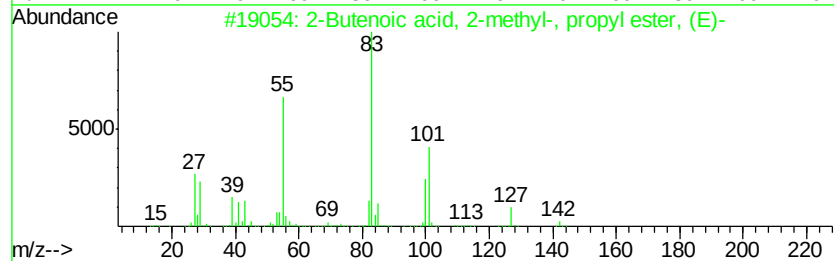
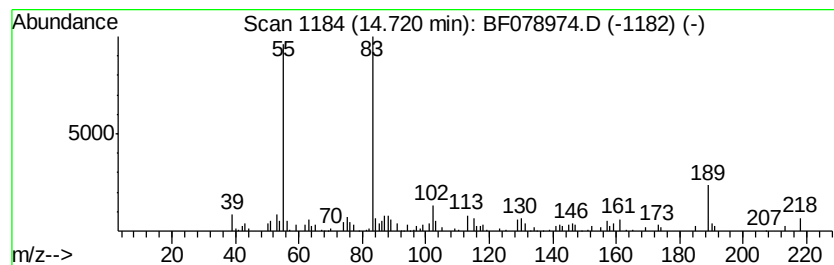
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Butenoic acid, 2-methyl-,... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	3.31 ng	167735	Chrysene-d12	16.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenoic acid, 2-methyl-, prop...	142	C8H14O2	061692-83-9	50
2		Cyclopropane, 2-chloro-1,1,3-tri...	118	C6H11Cl	098485-99-5	47
3		3-Hexen-2-one	98	C6H10O	000763-93-9	47
4		Cyclopentane acetic acid hydrazide	142	C7H14N2O	020287-25-6	47
5		Cyclobutanecarboxylic acid, 3,5-...	212	C11H10F2O2	1000293-70-4	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
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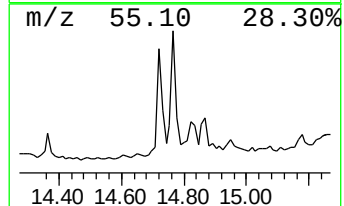
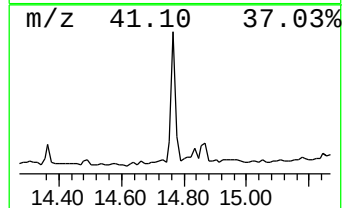
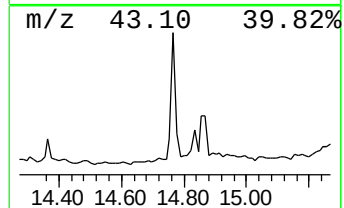
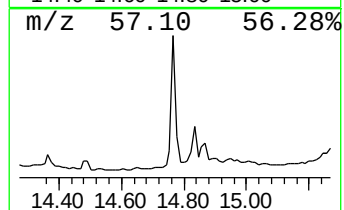
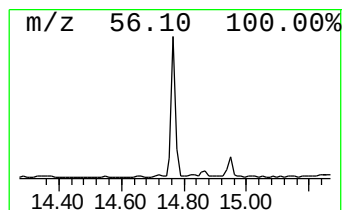
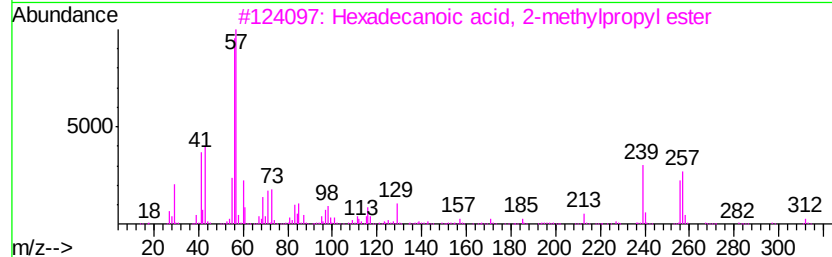
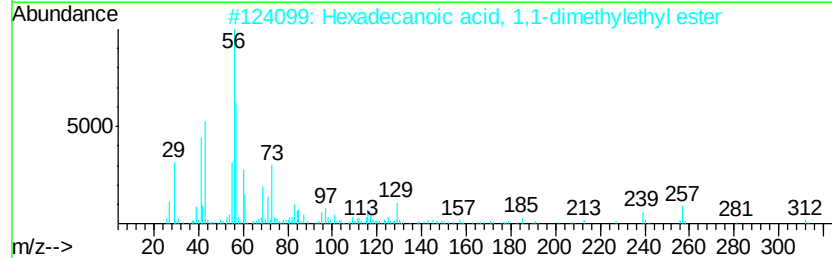
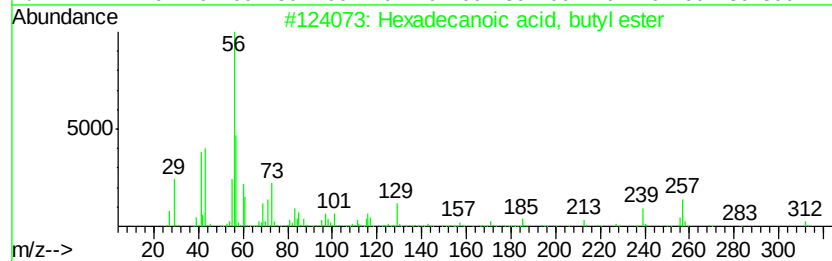
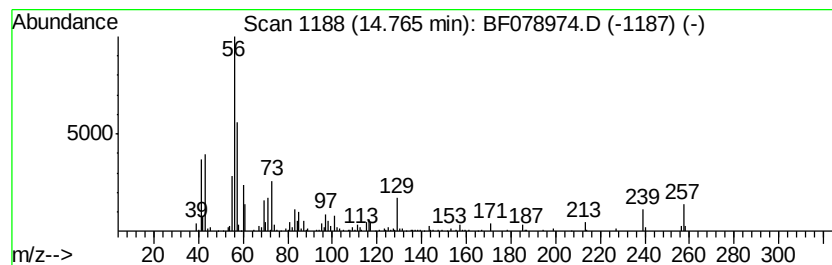
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecanoic acid, butyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	5.28 ng	267845	Chrysene-d12	16.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	91
3		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	43
4		Cyclohexanamine	99	C6H13N	000108-91-8	38
5		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
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Acq On : 6 May 2015 21:47
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Sample : G2035-17
Misc :
ALS Vial : 24 Sample Multiplier: 1

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ClientSampleId :
TP-15-SS

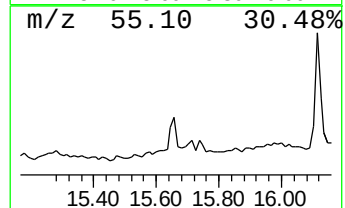
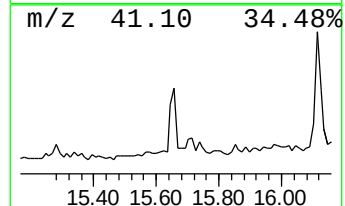
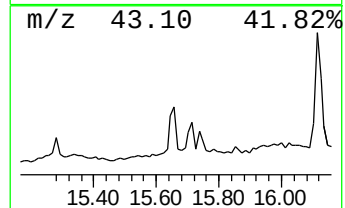
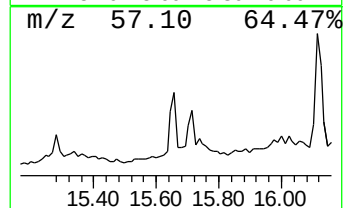
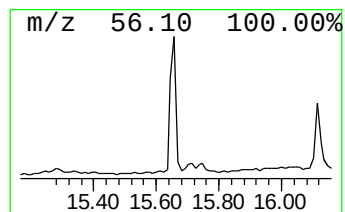
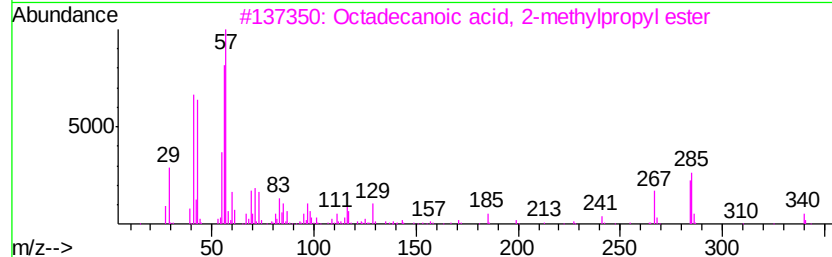
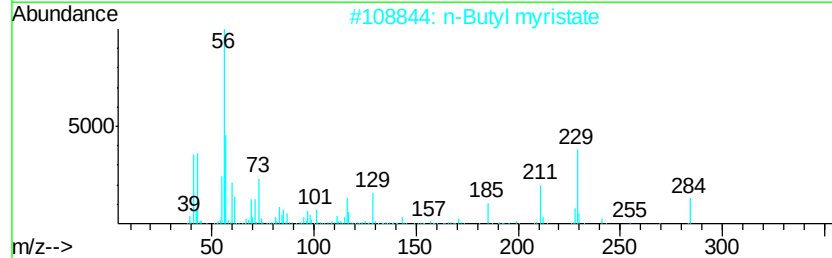
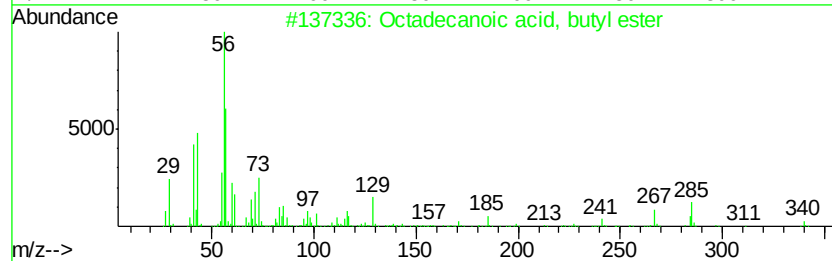
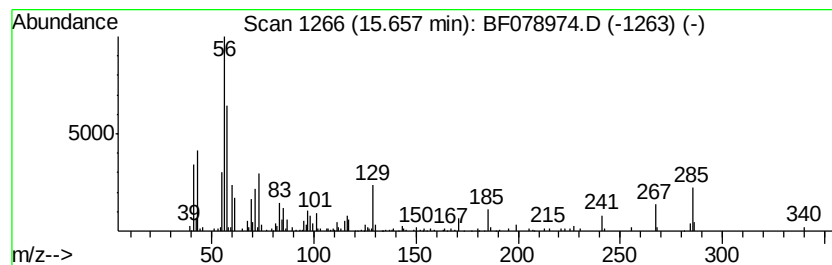
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid, butyl ester Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	4.23 ng	214466	Chrysene-d12	16.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	97
2		n-Butyl myristate	284	C18H36O2	000110-36-1	74
3		Octadecanoic acid, 2-methylpropyl ester	340	C22H44O2	000646-13-9	43
4		1-Hexene, 4-methyl-	98	C7H14	003769-23-1	30
5		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050615\
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Acq On : 6 May 2015 21:47
Operator : TP/IZ
Sample : G2035-17
Misc :
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Instrument :
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ClientSampleId :
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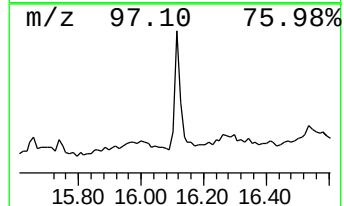
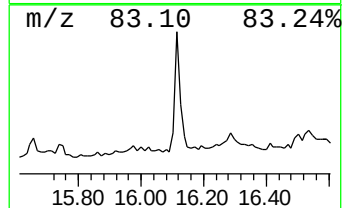
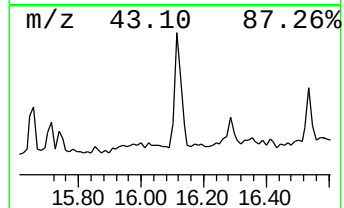
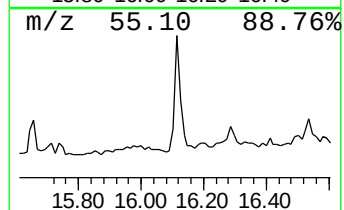
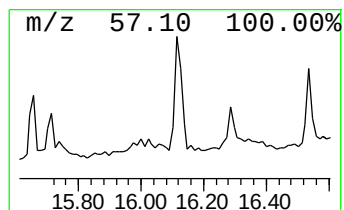
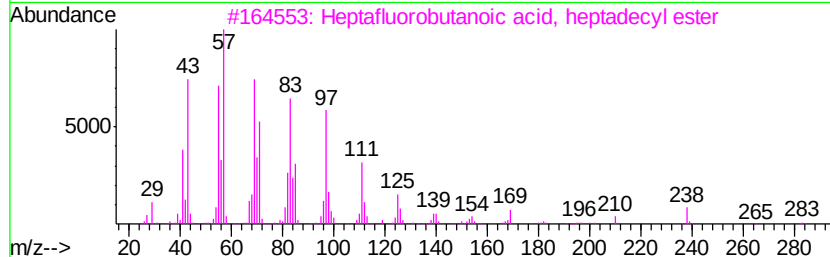
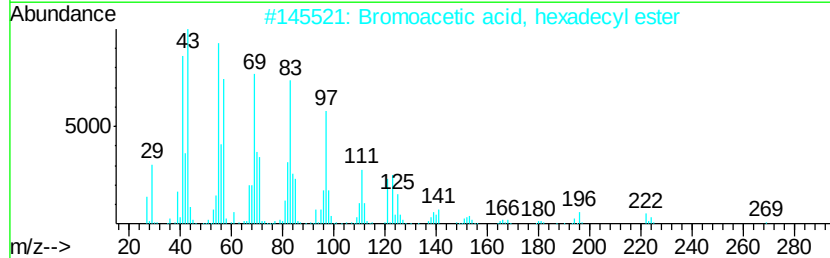
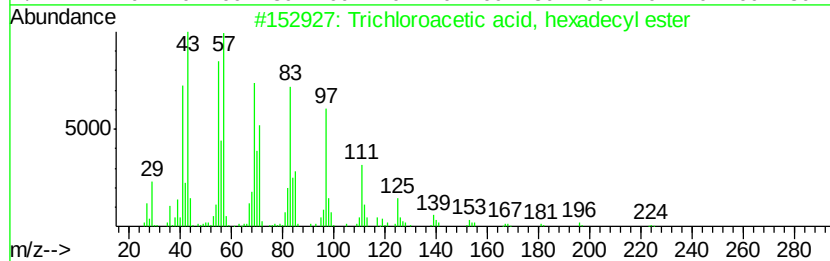
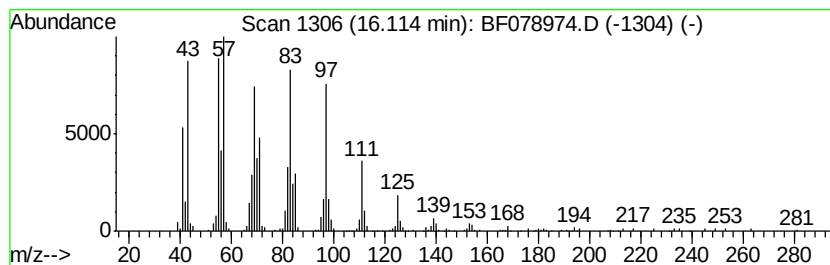
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Trichloroacetic acid, hexad... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.11	7.80 ng	395453	Chrysene-d12	16.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	91
3		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
5		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91



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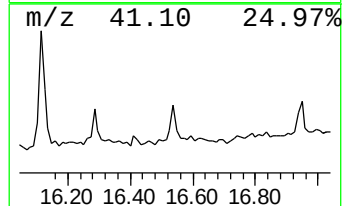
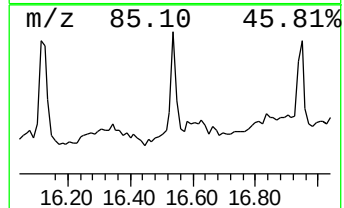
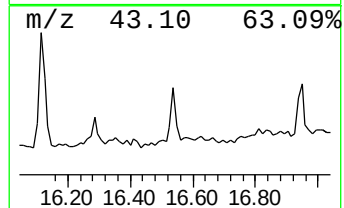
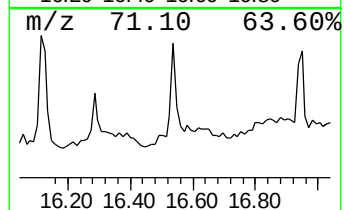
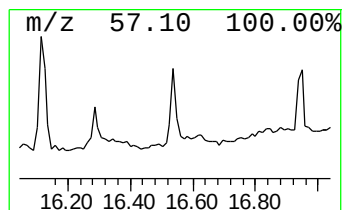
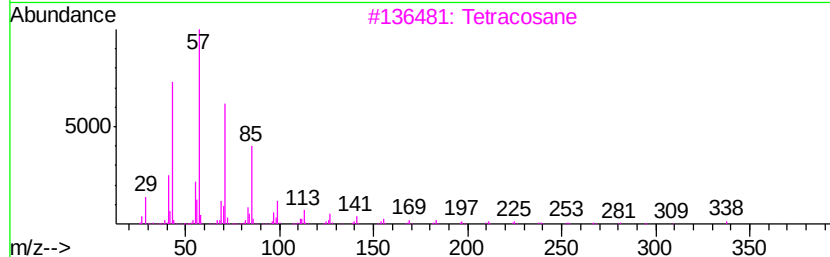
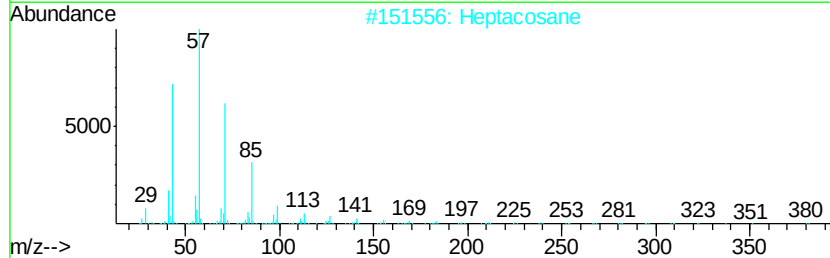
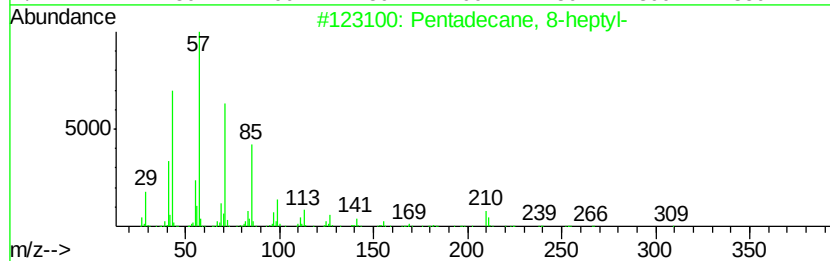
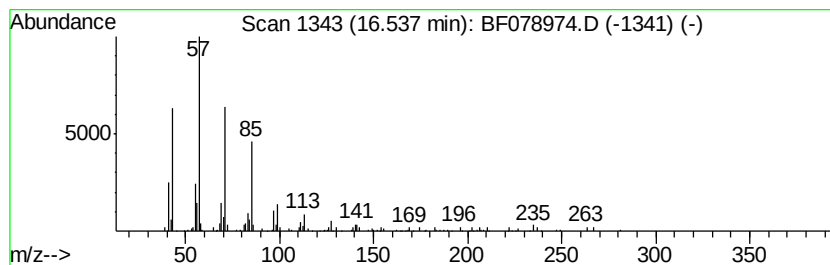
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pentadecane, 8-heptyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	2.64 ng	134011	Chrysene-d12	16.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	87
2		Heptacosane	380	C27H56	000593-49-7	83
3		Tetracosane	338	C24H50	000646-31-1	80
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	80
5		Nonadecane	268	C19H40	000629-92-5	80



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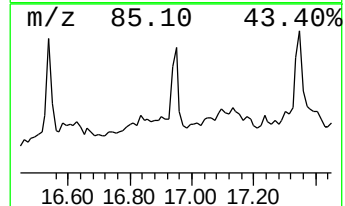
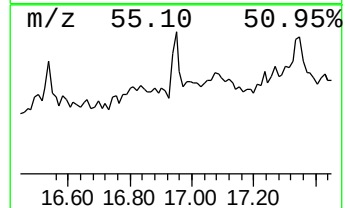
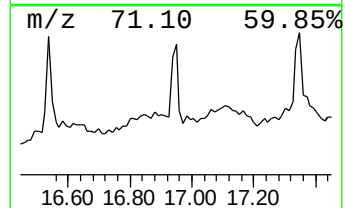
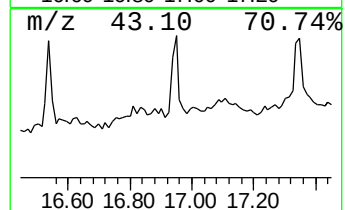
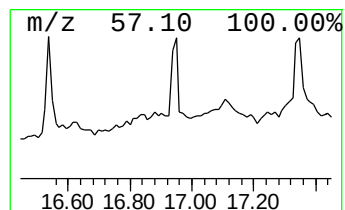
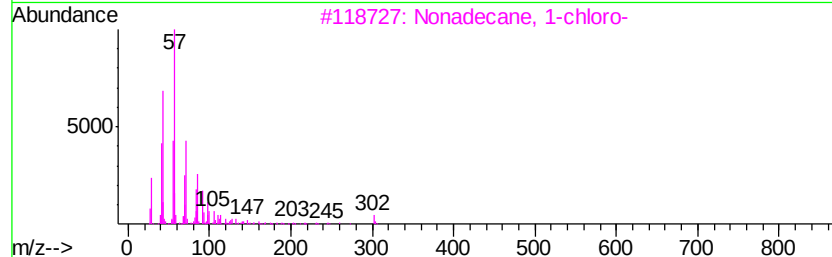
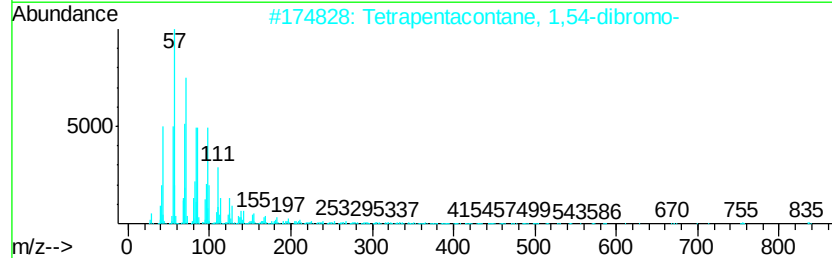
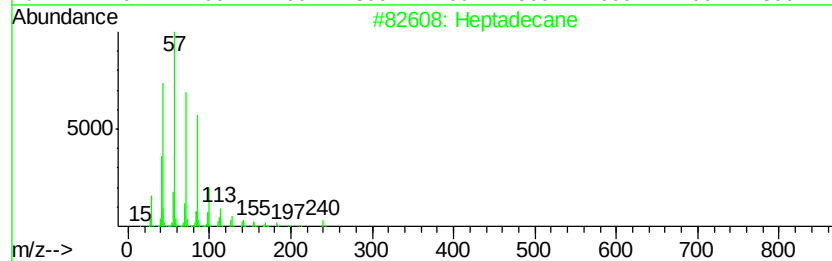
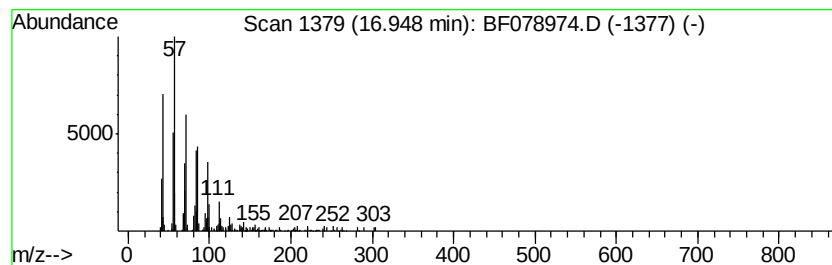
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heptadecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.95	3.29 ng	166621	Chrysene-d12	16.25
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	95
2	Tetrapentacontane, 1,54-dibromo-	915 C54H108Br2	1000156-09-4	90
3	Nonadecane, 1-chloro-	302 C19H39Cl	062016-76-6	76
4	Tetradecane	198 C14H30	000629-59-4	53
5	Hexadecane	226 C16H34	000544-76-3	53



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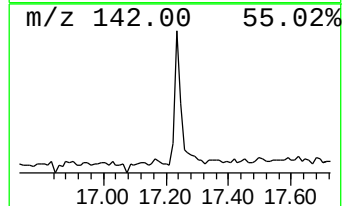
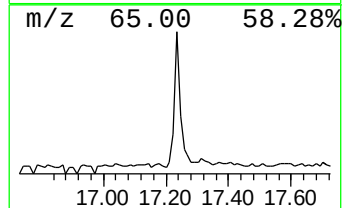
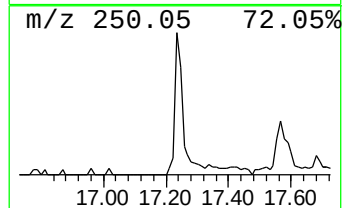
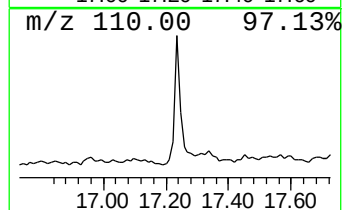
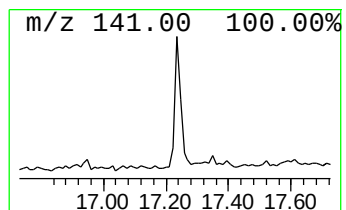
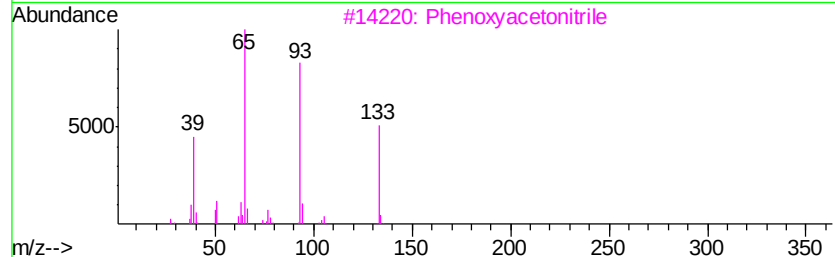
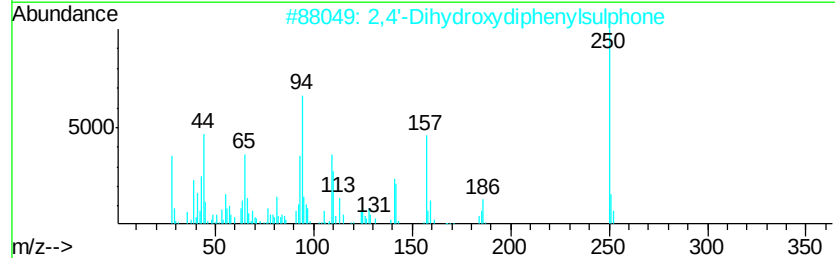
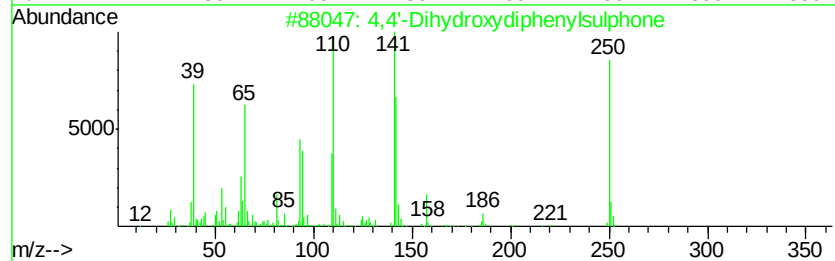
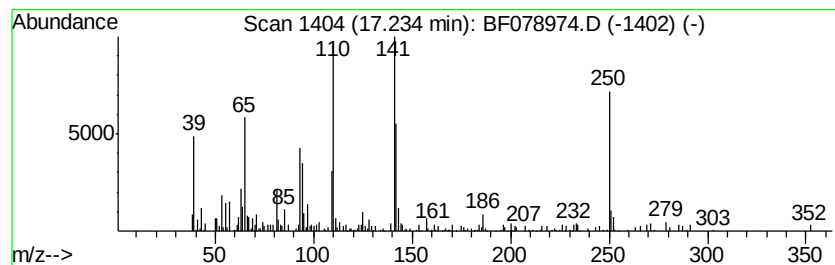
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 4,4'-Dihydroxydiphenylsulphone Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.23	3.56 ng	225451	Perylene-d12	18.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4,4'-Dihydroxydiphenylsulphone	250	C12H10O4S	000080-09-1	97
2		2,4'-Dihydroxydiphenylsulphone	250	C12H10O4S	005397-34-2	15
3		Phenoxvacetonitrile	133	C8H7NO	003598-14-9	12
4		1-Acetylcyclopentadiene	108	C7H8O	060032-12-4	12
5		n-Decyl(phenyl)sulfide	250	C16H26S	013910-18-4	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050615\
Data File : BF078974.D
Acq On : 6 May 2015 21:47
Operator : TP/IZ
Sample : G2035-17
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-15-SS

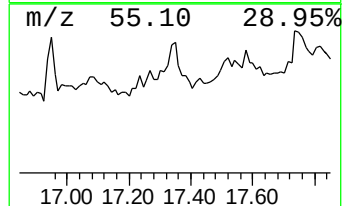
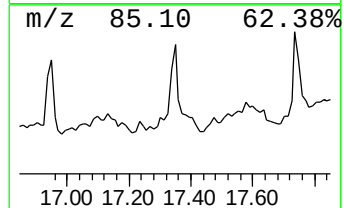
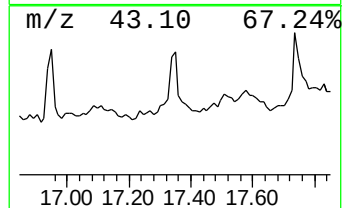
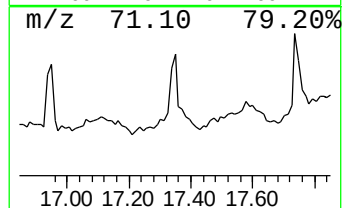
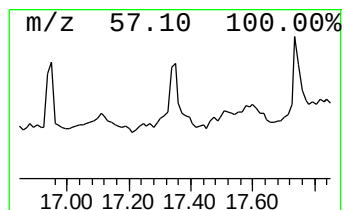
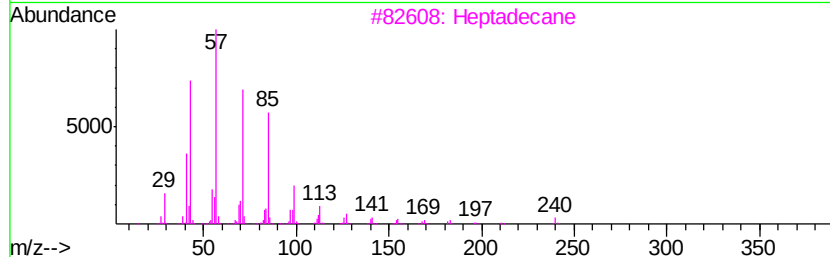
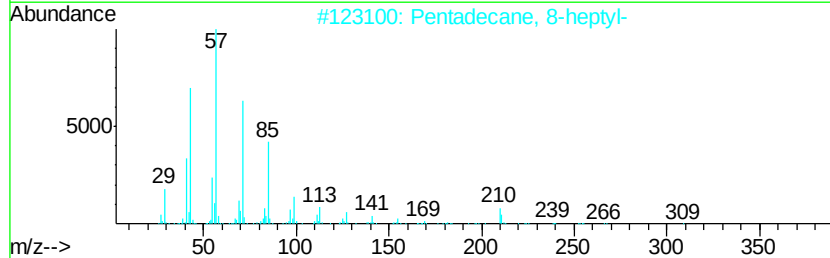
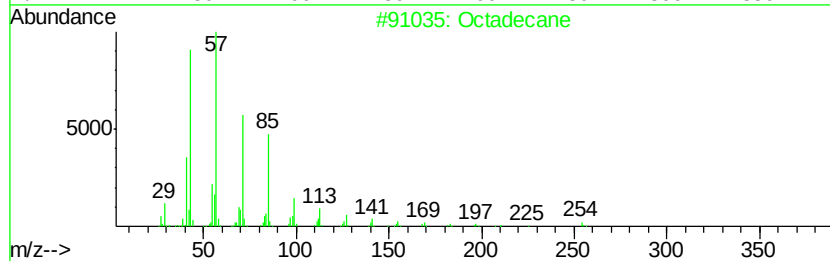
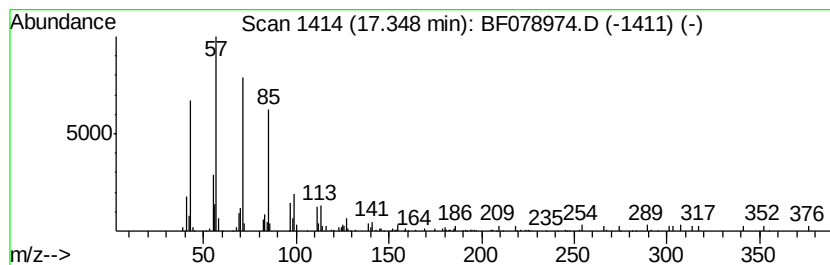
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Octadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.35	3.83 ng	242310	Perylene-d12	18.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	87
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
3		Heptadecane	240	C17H36	000629-78-7	83
4		Borane, diethyl(decyloxy)-	226	C14H31BO	1000152-34-3	76
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
Data File : BF078974.D
Acq On : 6 May 2015 21:47
Operator : TP/IZ
Sample : G2035-17
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-15-SS

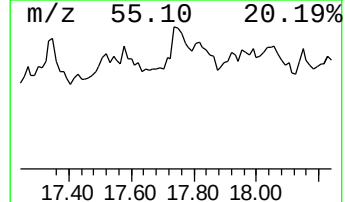
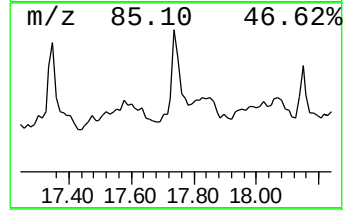
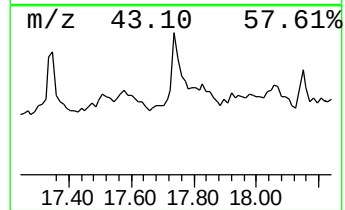
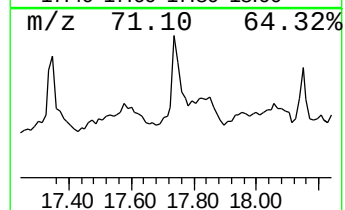
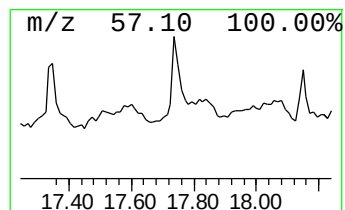
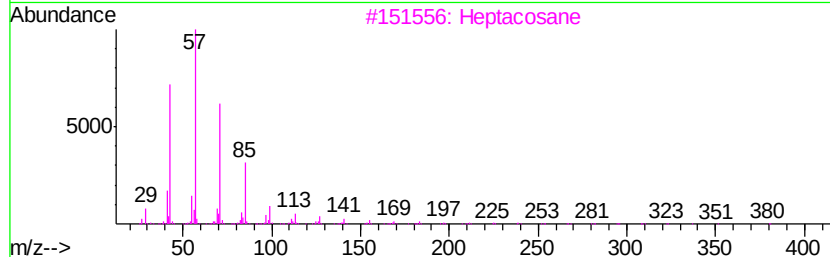
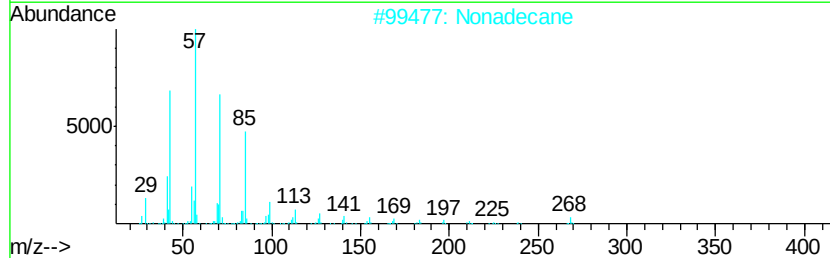
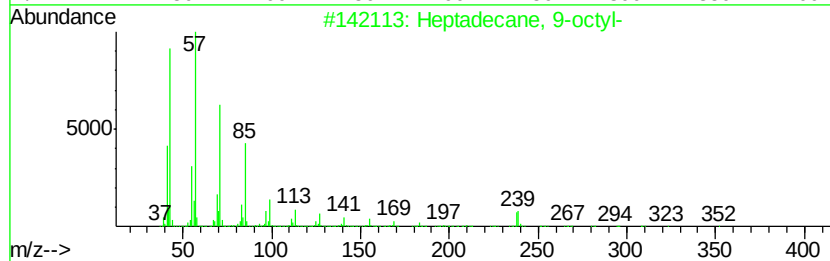
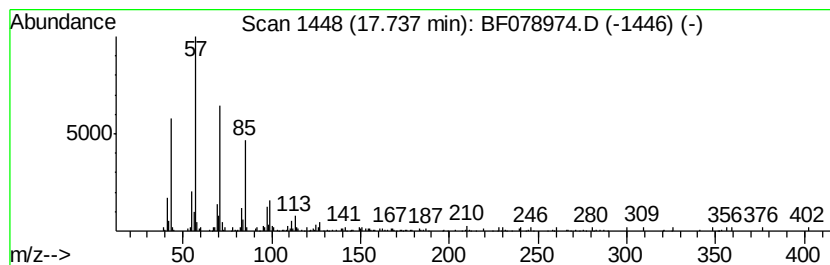
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Heptadecane, 9-octyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.74	3.45 ng	218344	Perylene-d12	18.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	86
2		Nonadecane	268	C19H40	000629-92-5	83
3		Heptacosane	380	C27H56	000593-49-7	80
4		Tetracosane	338	C24H50	000646-31-1	72
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
Data File : BF078974.D
Acq On : 6 May 2015 21:47
Operator : TP/IZ
Sample : G2035-17
Misc :
ALS Vial : 24 Sample Multiplier: 1

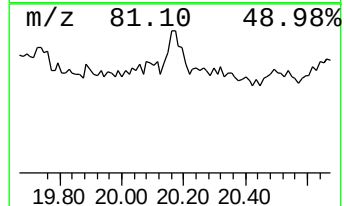
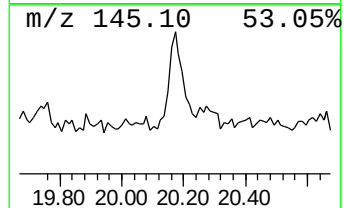
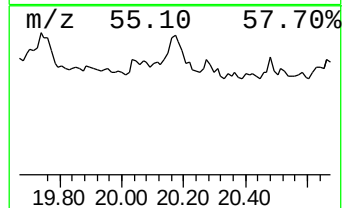
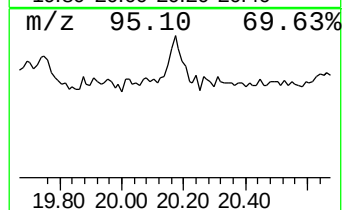
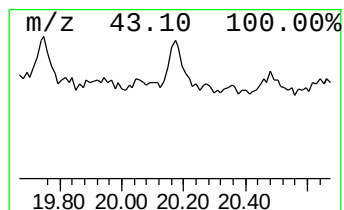
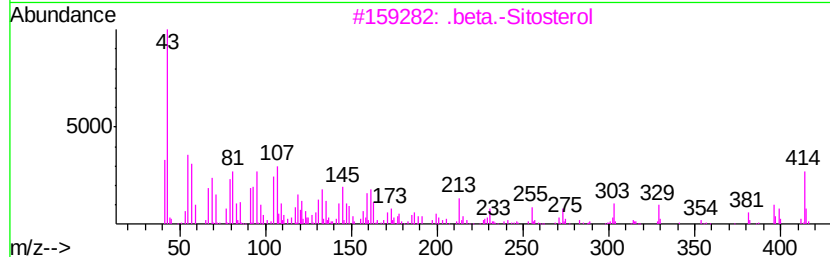
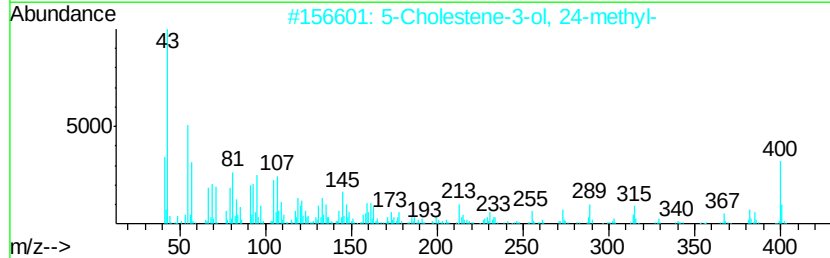
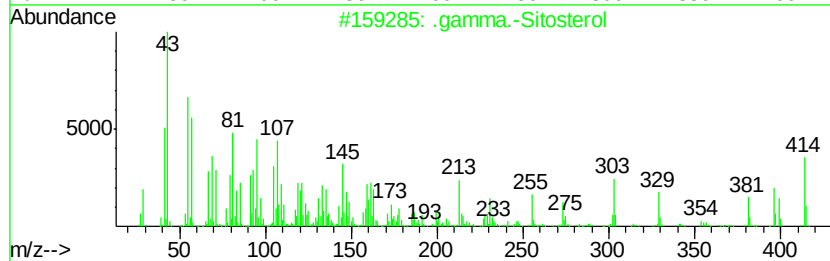
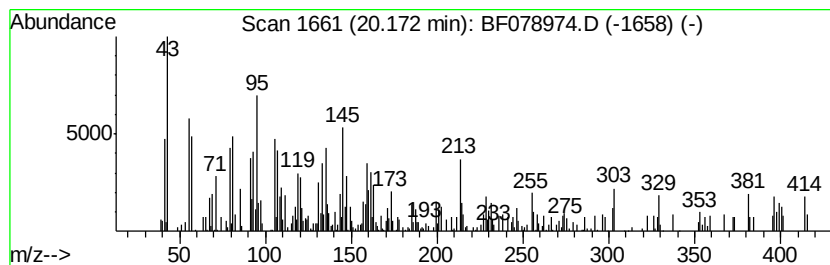
Instrument :
BNA_F
ClientSampleId :
TP-15-SS

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 20 .gamma.-Sitosterol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.17	4.79 ng	302688	Perylene-d12	18.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	86
2	5-Cholestene-3-ol, 24-methyl-	400	C28H48O	1000214-17-4	83
3	.beta.-Sitosterol	414	C29H50O	000083-46-5	64
4	Ergost-7-en-3-ol, (3.beta.)-	400	C28H48O	026047-31-4	38
5	Cholest-7-en-3-ol, 4,4-dimethyl-...	414	C29H50O	006384-28-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050615\
 Data File : BF078974.D
 Acq On : 6 May 2015 21:47
 Operator : TP/IZ
 Sample : G2035-17
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-15-SS

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
Propane, 2,2-dime...	1.50	334.3	ng	6937930	1	7.35	415138	20.0
Butane, 2-methoxy...	1.82	12.0	ng	248582	1	7.35	415138	20.0
2-Pentanone, 4-hy...	5.15	19.1	ng	396305	1	7.35	415138	20.0
unknown7.06	7.06	105.1	ng	2181750	1	7.35	415138	20.0
unknown12.03	12.03	5.2	ng	195740	4	12.95	758367	20.0
2-Butenoic acid, ...	14.72	3.3	ng	167735	5	16.25	1013930	20.0
Hexadecanoic acid...	14.77	5.3	ng	267845	5	16.25	1013930	20.0
Octadecanoic acid...	15.66	4.2	ng	214466	5	16.25	1013930	20.0
Trichloroacetic a...	16.11	7.8	ng	395453	5	16.25	1013930	20.0
Pentadecane, 8-he...	16.54	2.6	ng	134011	5	16.25	1013930	20.0
Heptadecane	16.95	3.3	ng	166621	5	16.25	1013930	20.0
4,4'-Dihydroxydip...	17.23	3.6	ng	225451	6	18.02	1265000	20.0
Octadecane	17.35	3.8	ng	242310	6	18.02	1265000	20.0
Heptadecane, 9-oc...	17.74	3.5	ng	218344	6	18.02	1265000	20.0
.gamma.-Sitosterol	20.17	4.8	ng	302688	6	18.02	1265000	20.0