

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011915\
 Data File : BF076935.D
 Acq On : 19 Jan 2015 18:16
 Operator : IZ / TP
 Sample : G1101-24
 Misc : W
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB-2

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-BF011415.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	1.452	31	32	40	rBV	29574	67400	4.98%	0.722%
2	4.116	263	265	270	rBV	15451	33023	2.44%	0.354%
3	5.030	343	345	354	rBV	189463	365115	26.99%	3.911%
4	7.362	547	549	555	rBV	106444	203127	15.02%	2.176%
5	7.465	555	558	566	rVB	530679	819002	60.54%	8.773%
6	7.967	599	602	607	rBV	112743	192393	14.22%	2.061%
7	8.299	627	631	634	rBV	472436	813883	60.17%	8.718%
8	8.962	686	689	695	rBV2	8295	14174	1.05%	0.152%
9	9.270	712	716	719	rBV	436292	645385	47.71%	6.913%
10	10.882	854	857	861	rBV	171520	261835	19.36%	2.805%
11	11.396	899	902	904	rVB	25476	38250	2.83%	0.410%
12	12.139	965	967	972	rBV	29802	48435	3.58%	0.519%
13	13.534	1085	1089	1092	rBV	886035	1288097	95.22%	13.798%
14	13.957	1123	1126	1130	rBV	166501	267261	19.76%	2.863%
15	14.757	1194	1196	1201	rBV	40527	62563	4.62%	0.670%
16	15.008	1215	1218	1222	rVB	191422	285686	21.12%	3.060%
17	15.260	1236	1240	1243	rVB	72470	114161	8.44%	1.223%
18	15.362	1246	1249	1252	rVB	46268	68022	5.03%	0.729%
19	16.311	1329	1332	1334	rVB	39868	51457	3.80%	0.551%
20	16.471	1342	1346	1349	rBV	321691	471534	34.86%	5.051%
21	16.677	1361	1364	1367	rBV	570795	762459	56.36%	8.168%
22	17.774	1457	1460	1463	rBV	251865	289731	21.42%	3.104%
23	18.460	1517	1520	1522	rBV	19966	20572	1.52%	0.220%
24	19.900	1644	1646	1649	rVB	110111	107237	7.93%	1.149%
25	20.003	1652	1655	1658	rVB	1281070	1352746	100.00%	14.491%
26	20.792	1722	1724	1727	rVB	75913	73887	5.46%	0.791%
27	20.929	1734	1736	1740	rBV	10100	14320	1.06%	0.153%
28	21.272	1763	1766	1769	rBV	293874	309690	22.89%	3.317%
29	22.940	1909	1912	1915	rBV	233443	273458	20.22%	2.929%
30	24.541	2048	2052	2055	rBV4	10174	20245	1.50%	0.217%

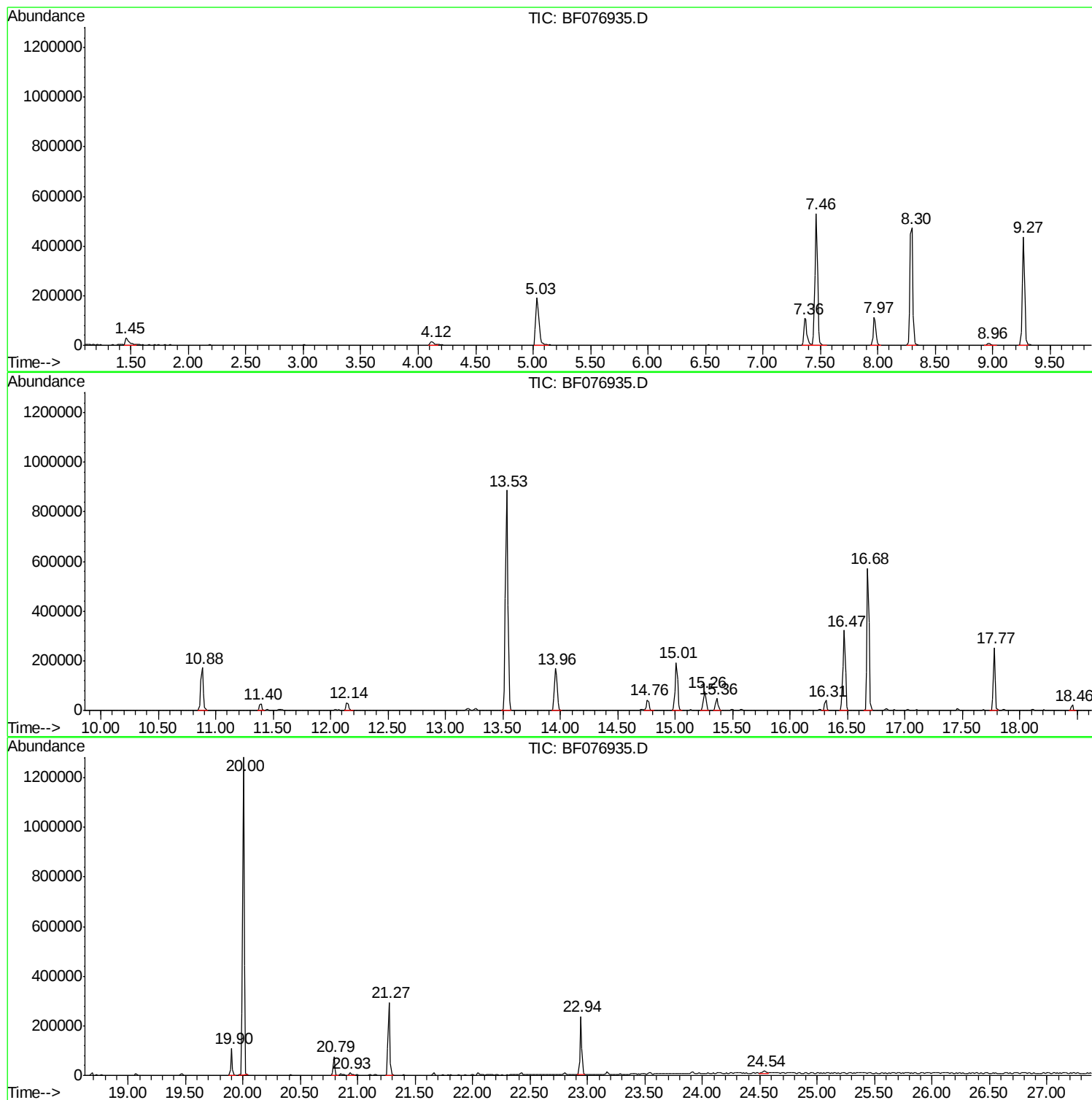
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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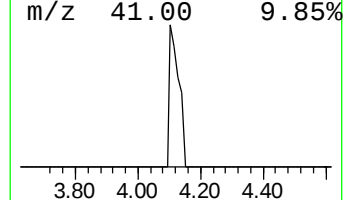
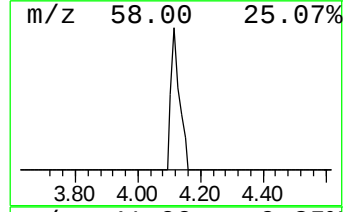
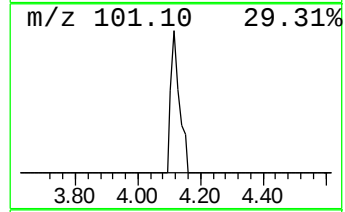
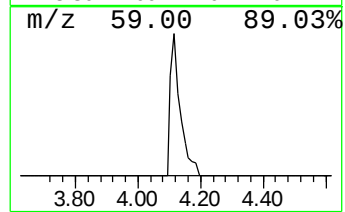
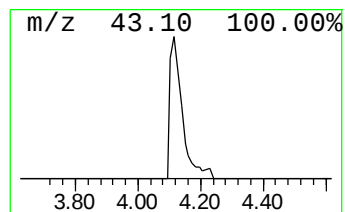
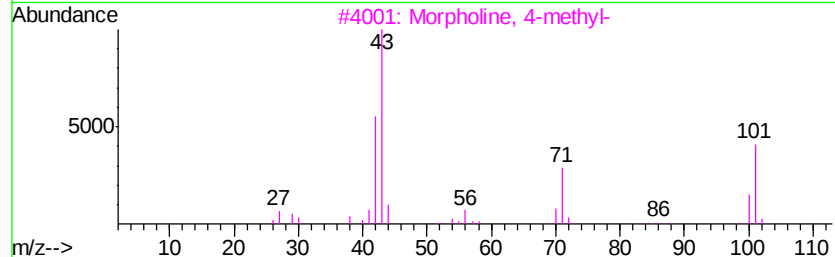
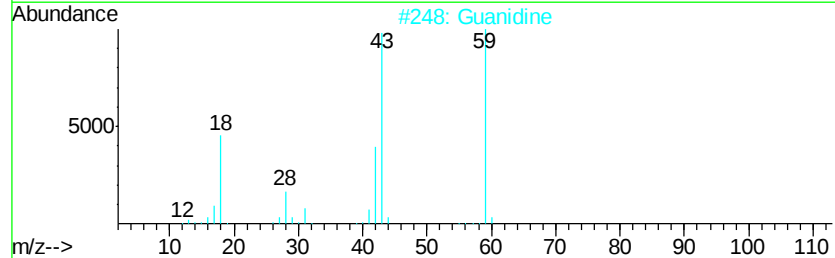
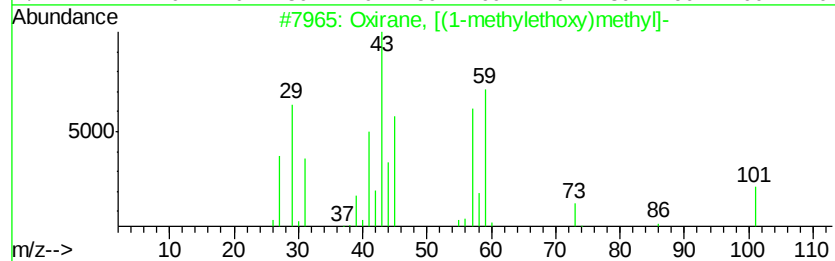
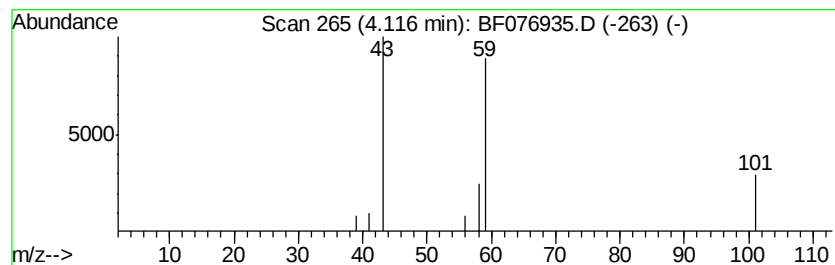
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Oxirane, [(1-methylethoxy)m... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.12	3.43 ng	33023	1,4-Dichlorobenzene-d4	7.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, [(1-methylethoxy)methyl]-	116	C6H12O2	004016-14-2	64
2			Guanidine	59	CH5N3	000113-00-8	7
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	5
4			Diacetamide	101	C4H7NO2	000625-77-4	5
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	5



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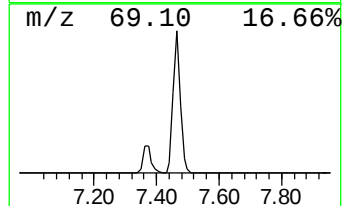
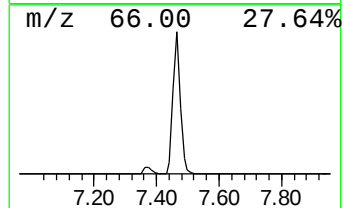
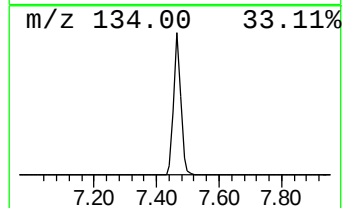
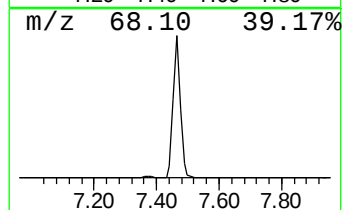
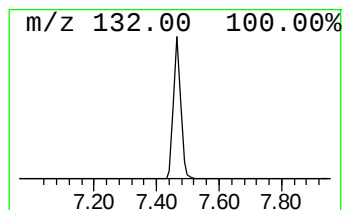
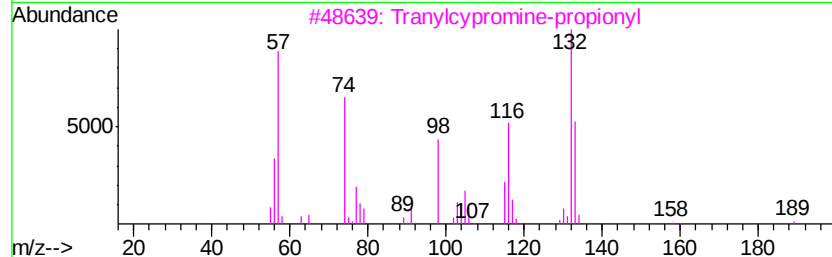
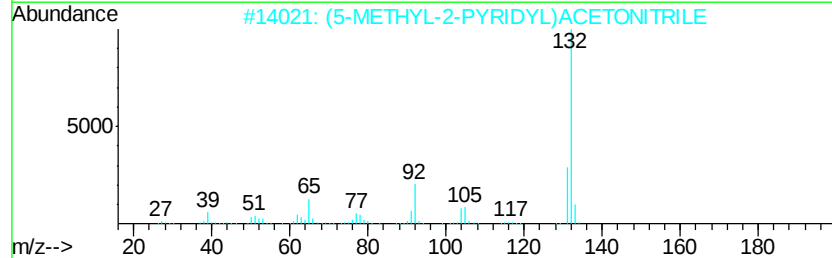
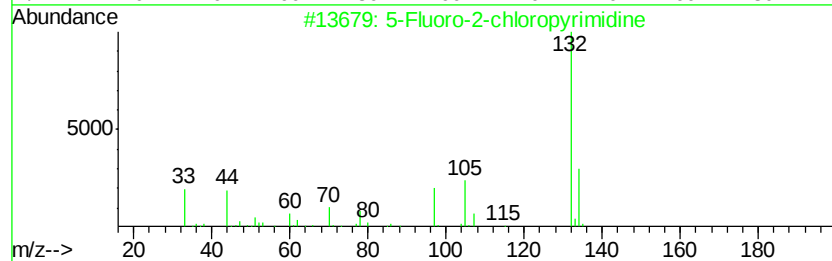
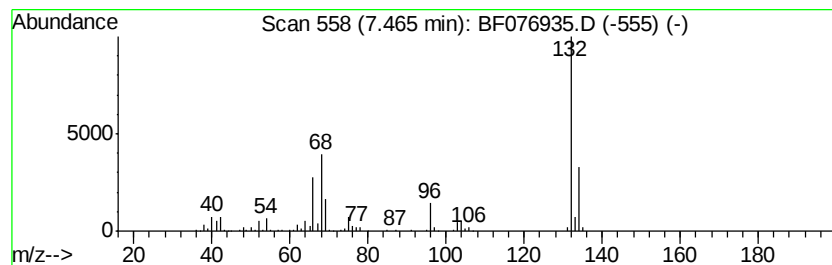
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Peak Number 3 unknown7.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.46	85.14 ng	819002	1,4-Dichlorobenzene-d4	7.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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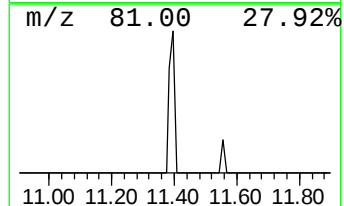
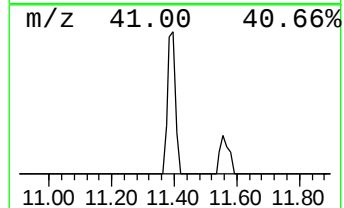
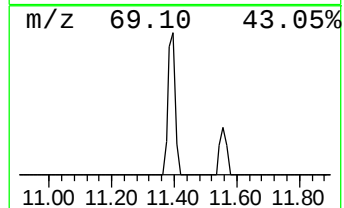
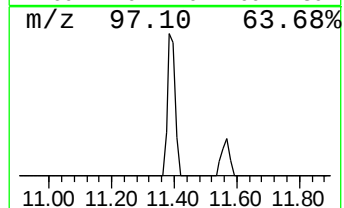
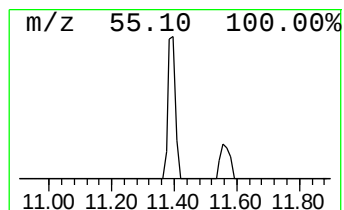
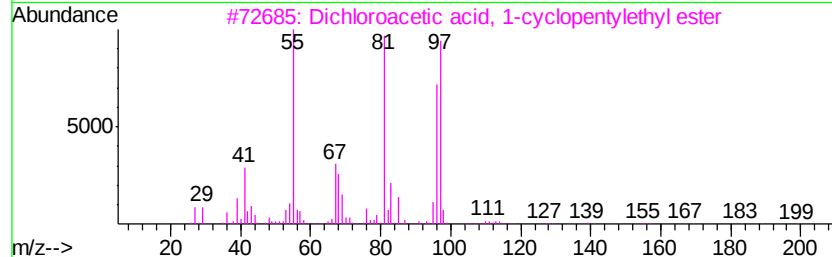
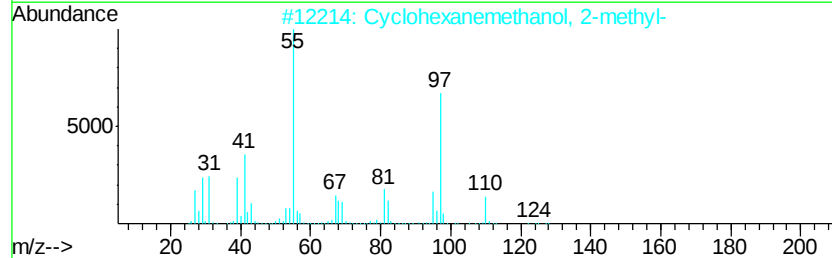
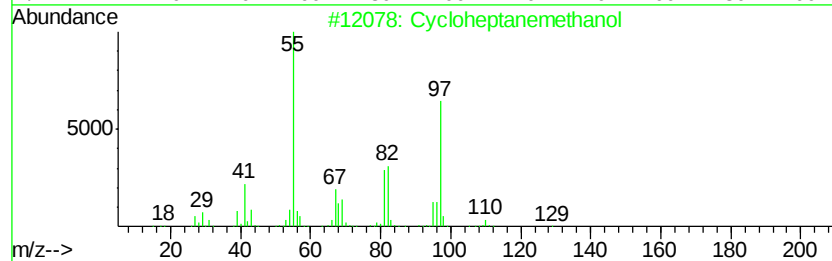
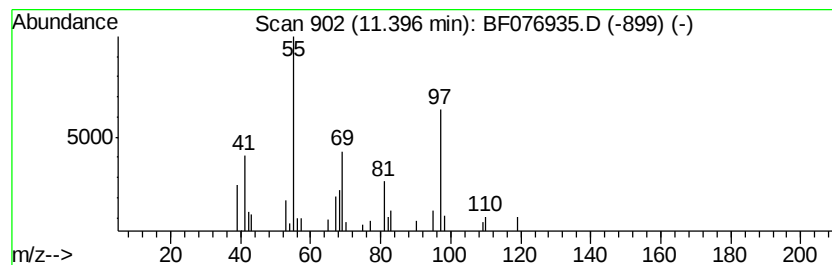
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Peak Number 5 Cycloheptanemethanol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	2.92 ng	38250	Naphthalene-d8	10.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cycloheptanemethanol	128	C8H16O	004448-75-3	53
2		Cyclohexanemethanol, 2-methyl-	128	C8H16O	002105-40-0	42
3		Dichloroacetic acid, 1-cyclopent...	224	C9H14Cl2O2	1000282-56-3	42
4		Cyclohexane, 1-bromo-2-methyl-	176	C7H13Br	006294-39-9	40
5		3-Hexen-1-ol, 2-ethyl-	128	C8H16O	053907-73-6	40



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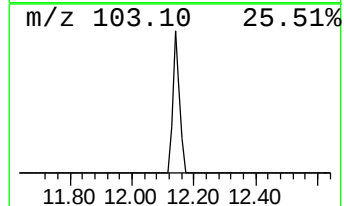
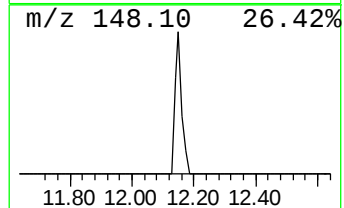
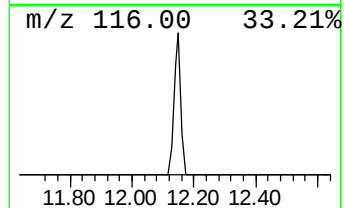
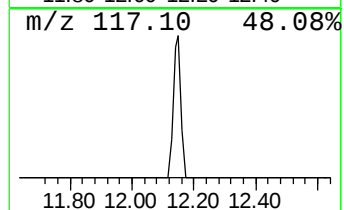
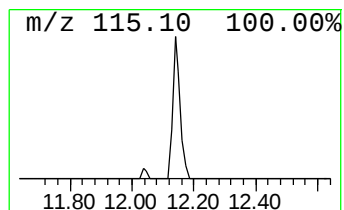
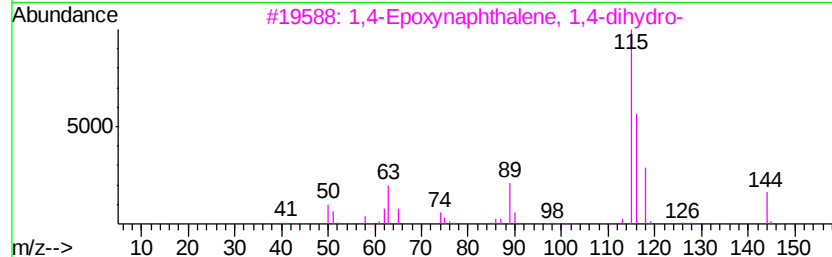
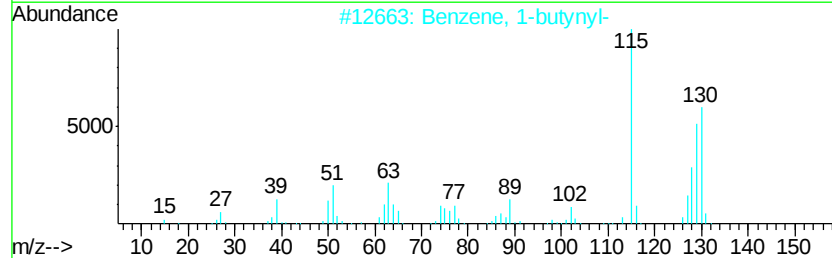
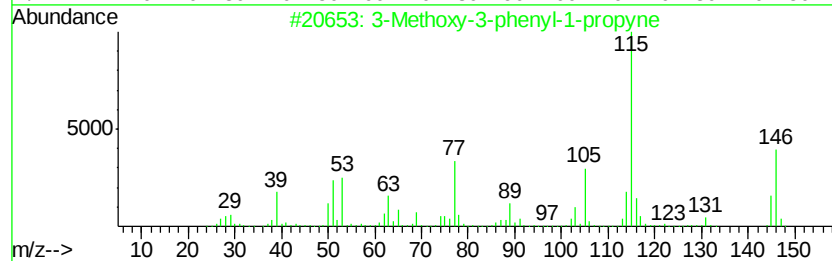
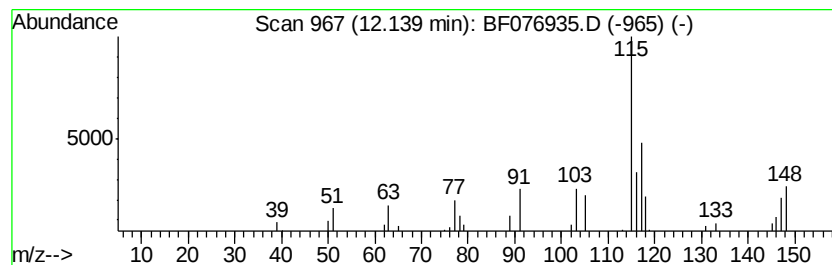
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Peak Number 6 unknown12.14 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	3.70 ng	48435	Naphthalene-d8	10.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methoxy-3-phenyl-1-propyne	146	C10H10O	1000217-10-1	43
2			Benzene, 1-butynyl-	130	C10H10	000622-76-4	22
3			1,4-Epoxy-naphthalene, 1,4-dihydro-	144	C10H8O	000573-57-9	10
4			Benzene, 1-ethynyl-2-methyl-	116	C9H8	000766-47-2	9
5			4-Phenylbut-3-yn-1-ol	146	C10H10O	010229-11-5	9



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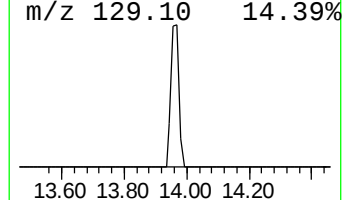
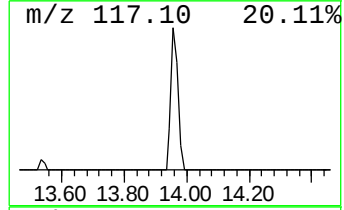
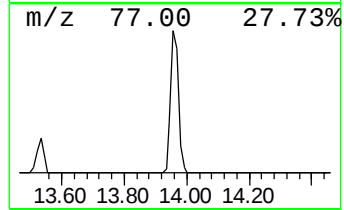
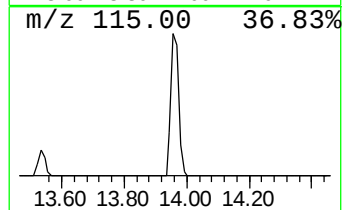
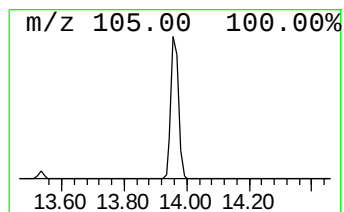
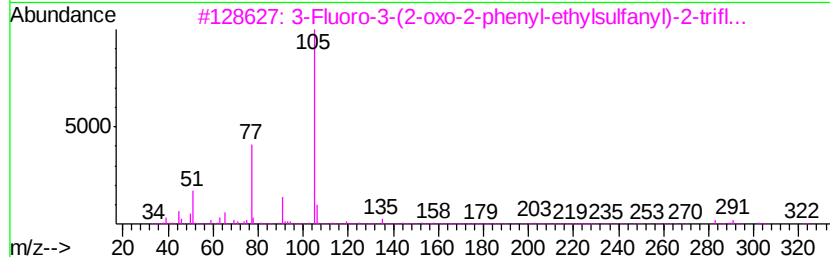
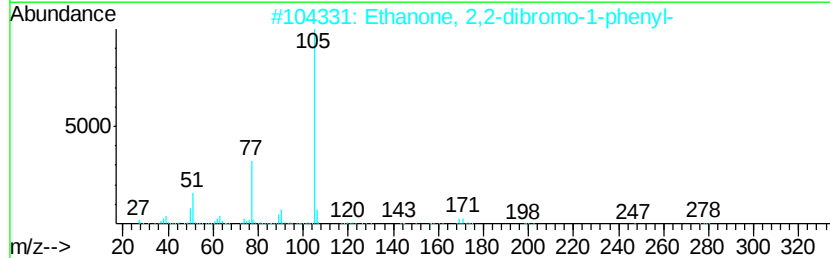
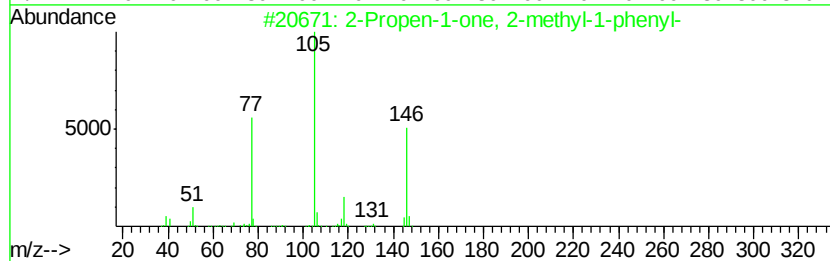
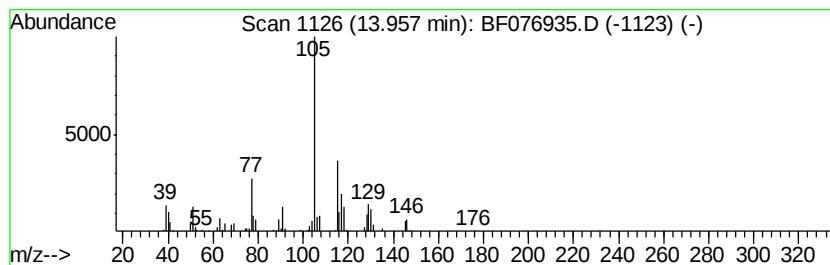
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown13.96 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	18.71 ng	267261	Acenaphthene-d10	15.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propen-1-one, 2-methyl-1-phenyl-	146	C10H10O	000769-60-8	35
2		Ethanone, 2,2-dibromo-1-phenyl-	276	C8H6Br2O	013665-04-8	14
3		3-Fluoro-3-(2-oxo-2-phenyl-ethyl)-...	322	C13H10F4O3S	329735-34-4	14
4		1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	14
5		Benzoic acid, 3,3,5-trimethyl-6-...	336	C21H20O4	1000277-63-9	14



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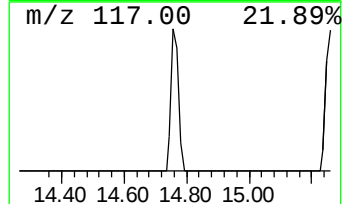
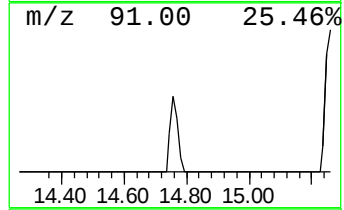
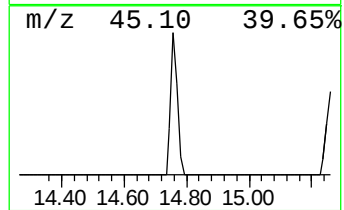
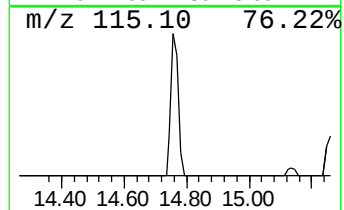
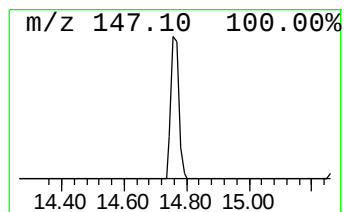
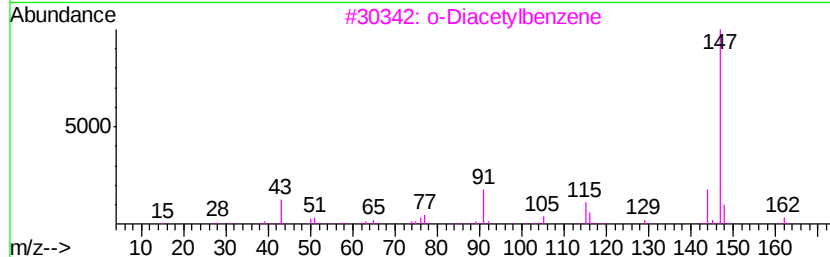
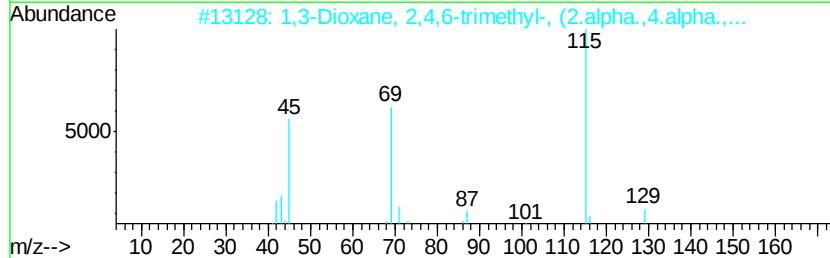
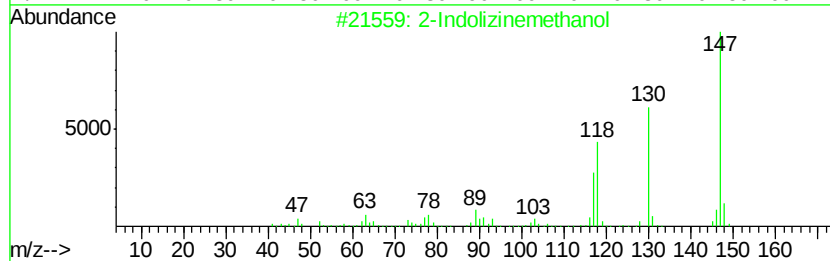
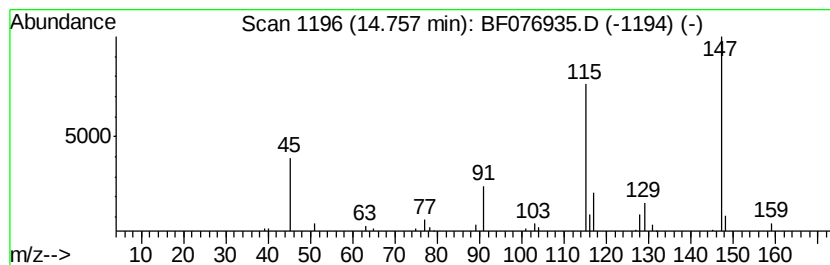
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.76 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	4.38 ng	62563	Acenaphthene-d10	15.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Indolizinemethanol	147	C9H9NO	022320-21-4	35
2		1,3-Dioxane, 2,4,6-trimethyl-, (...	130	C7H14O2	019145-91-6	23
3		o-Diacetylbenzene	162	C10H10O2	000704-00-7	16
4		2-(5-Methyl-[1,3,4]thiadiazol-2-...	292	C14H16N2OS2	1000296-87-8	16
5		2,5-Dimethylbenzoxazole	147	C9H9NO	005676-58-4	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011915\
Data File : BF076935.D
Acq On : 19 Jan 2015 18:16
Operator : IZ / TP
Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FB-2

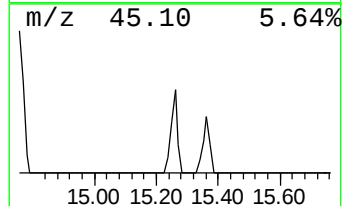
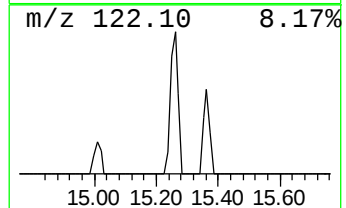
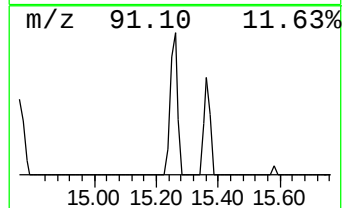
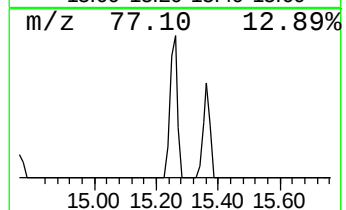
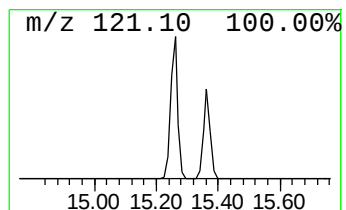
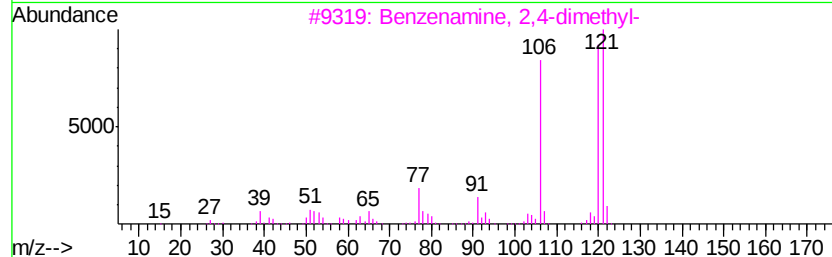
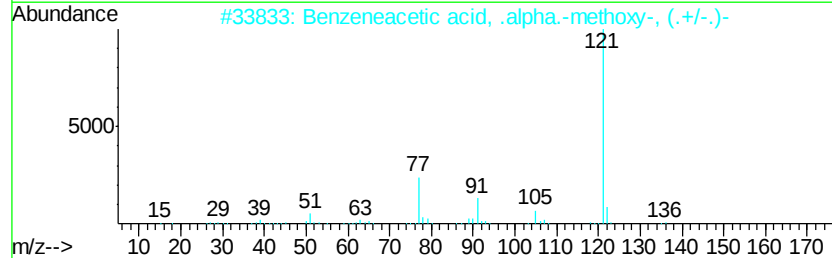
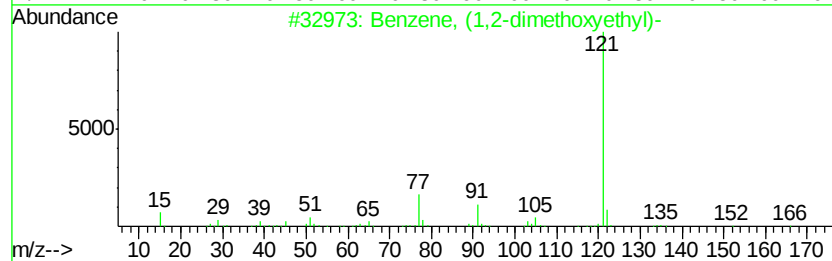
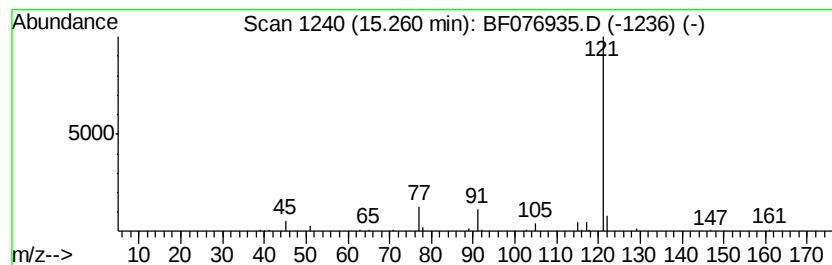
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, (1,2-dimethoxyethyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	7.99 ng	114161	Acenaphthene-d10	15.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1,2-dimethoxyethyl)-	166	C10H14O2	004013-37-0	72
2		Benzeneacetic acid, .alpha.-meth...	166	C9H10O3	007021-09-2	56
3		Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	40
4		dl-2-Phenyl-1,2-propanediol	152	C9H12O2	004217-66-7	40
5		Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011915\
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Acq On : 19 Jan 2015 18:16
Operator : IZ / TP
Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FB-2

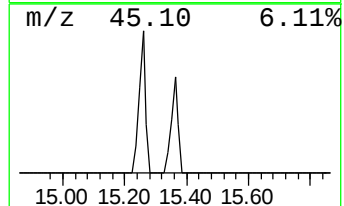
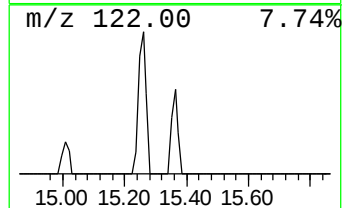
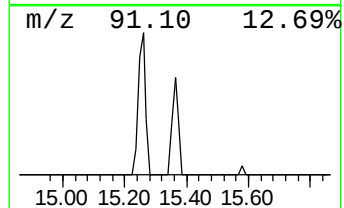
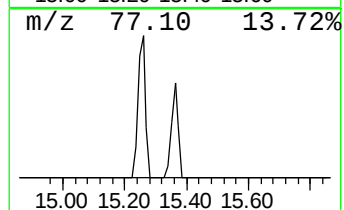
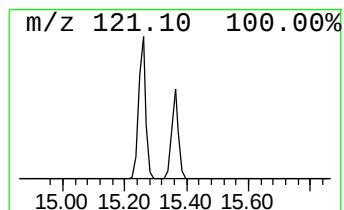
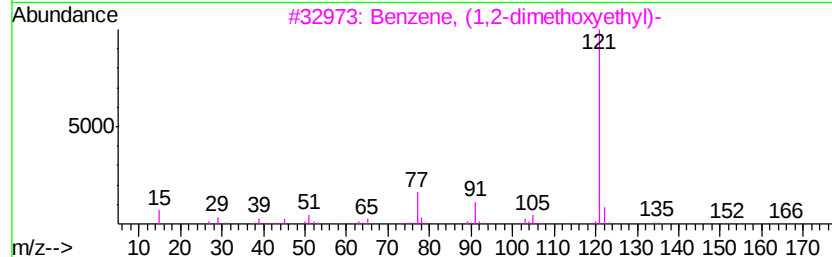
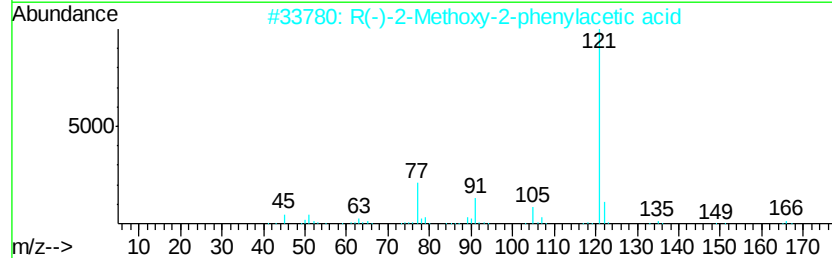
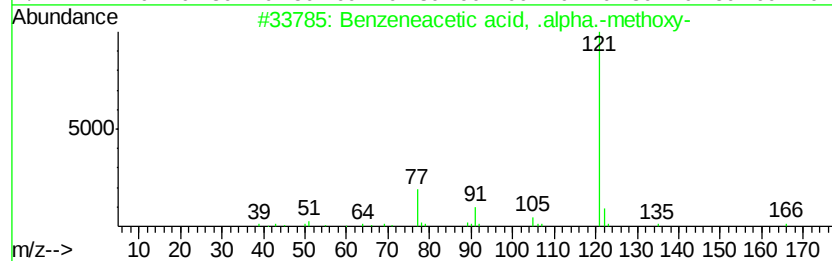
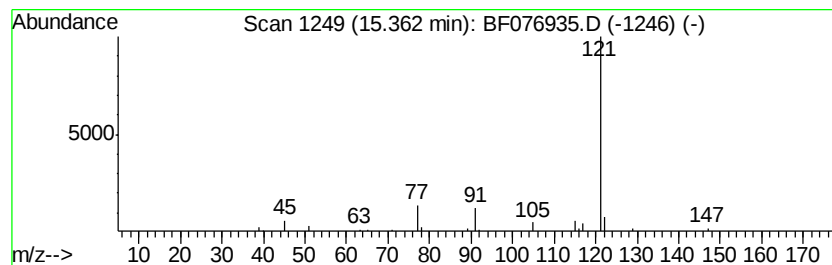
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzeneacetic acid, .alpha.... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	4.76 ng	68022	Acenaphthene-d10	15.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetic acid, .alpha.-meth...	166	C9H10O3	001701-77-5	72
2		R(-)-2-Methoxy-2-phenylacetic acid	166	C9H10O3	1000132-09-7	72
3		Benzene, (1,2-dimethoxyethyl)-	166	C10H14O2	004013-37-0	64
4		Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	56
5		Benzenamine, 3,4-dimethyl-	121	C8H11N	000095-64-7	56



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011915\
Data File : BF076935.D
Acq On : 19 Jan 2015 18:16
Operator : IZ / TP
Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

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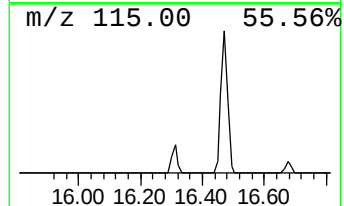
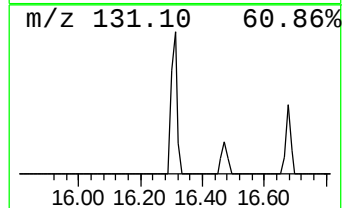
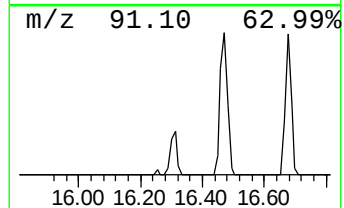
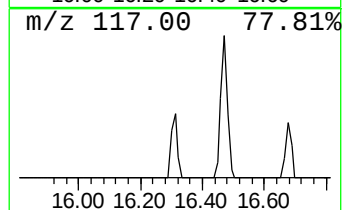
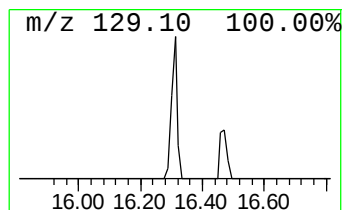
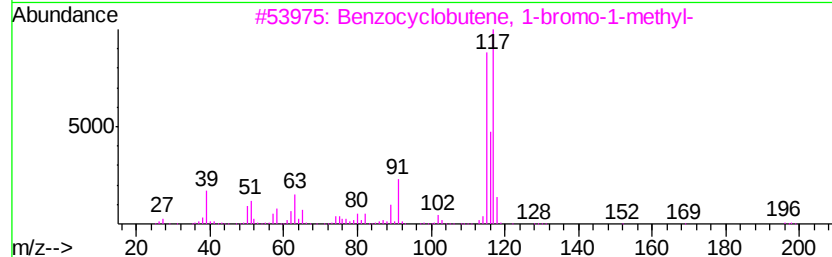
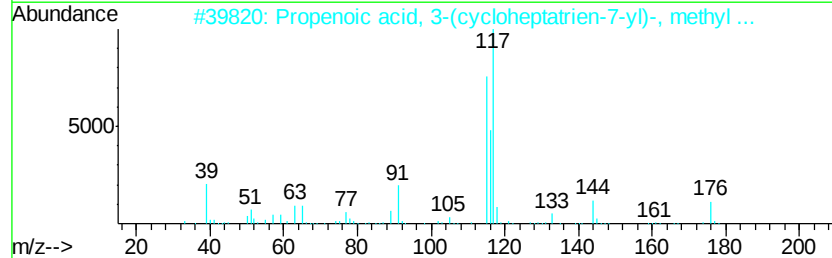
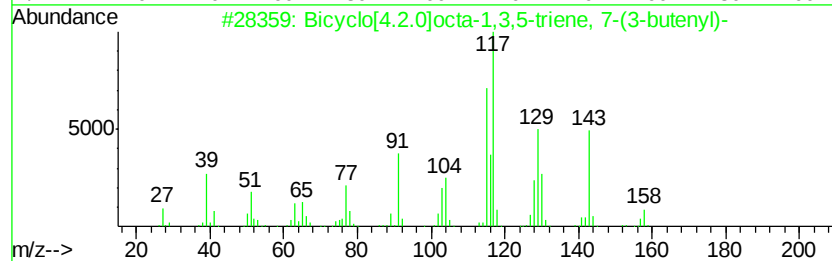
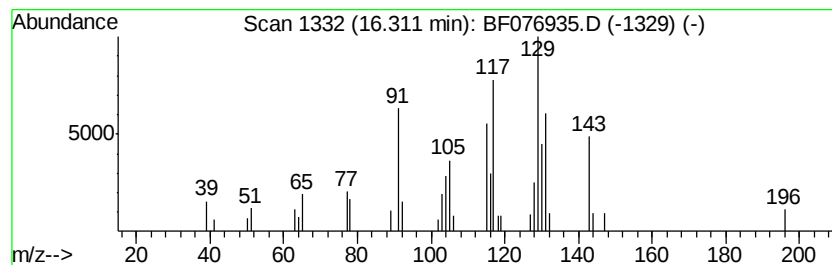
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.31	3.60 ng	51457	Acenaphthene-d10	15.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.2.0]octa-1,3,5-triene,...	158	C12H14	122057-61-8	58
2		Propenoic acid, 3-(cycloheptatrien...	176	C11H12O2	1000269-64-5	22
3		Benzocyclobutene, 1-bromo-1-methyl-	196	C9H9Br	1000161-76-4	22
4		Triquinacene	130	C10H10	006053-74-3	14
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011915\
Data File : BF076935.D
Acq On : 19 Jan 2015 18:16
Operator : IZ / TP
Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FB-2

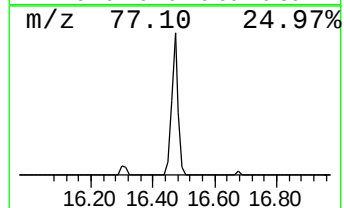
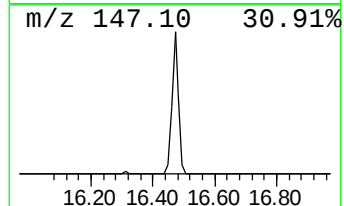
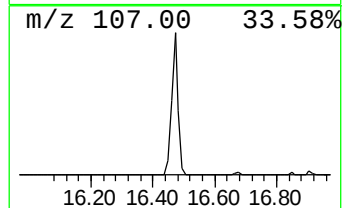
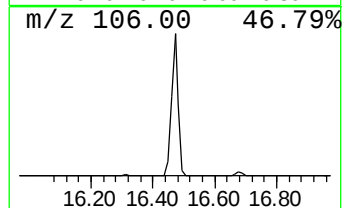
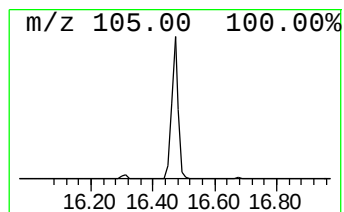
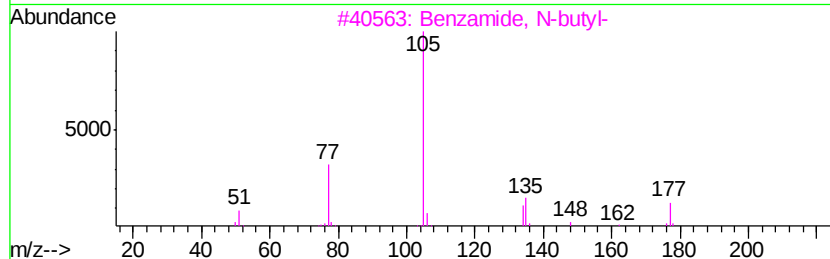
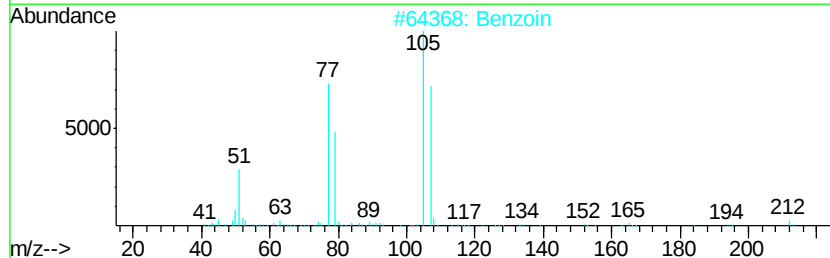
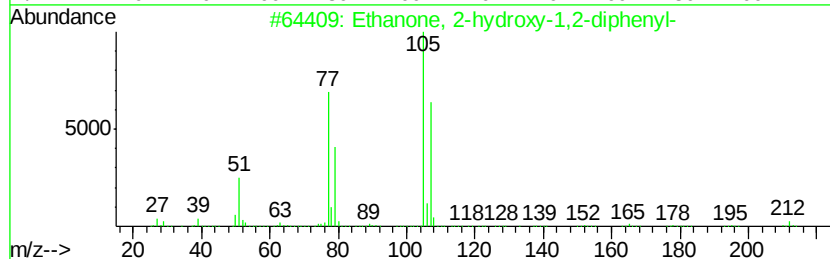
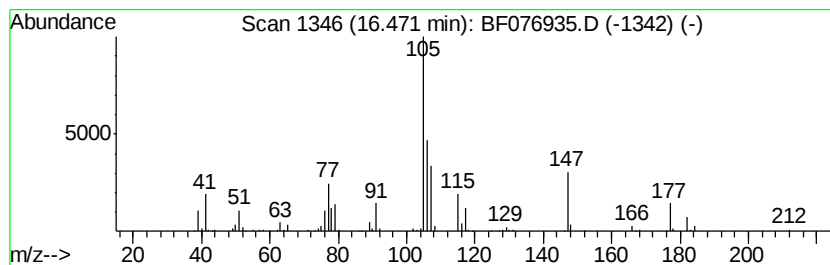
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown16.47 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	32.55 ng	471534	Phenanthrene-d10	17.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 2-hydroxy-1,2-diphenyl-	212	C14H12O2	000119-53-9	27
2		Benzoin	212	C14H12O2	000579-44-2	25
3		Benzamide, N-butyl-	177	C11H15NO	002782-40-3	25
4		4-Methylphenethyl isothiocyanate	177	C10H11NS	013203-39-9	22
5		Ethanone, 2-(acetyloxy)-1,2-diph...	254	C16H14O3	000574-06-1	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011915\
Data File : BF076935.D
Acq On : 19 Jan 2015 18:16
Operator : IZ / TP
Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FB-2

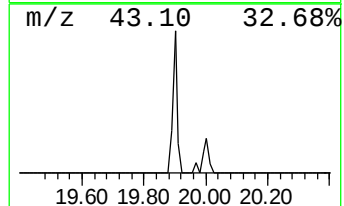
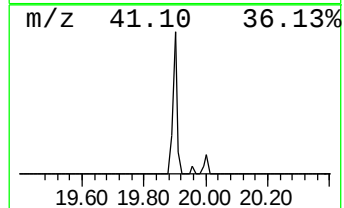
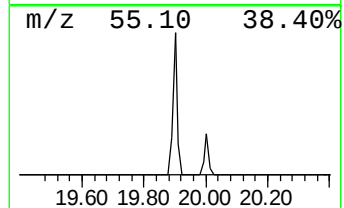
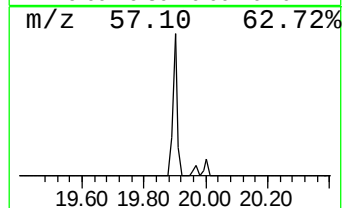
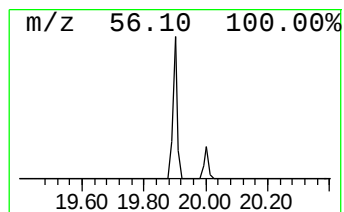
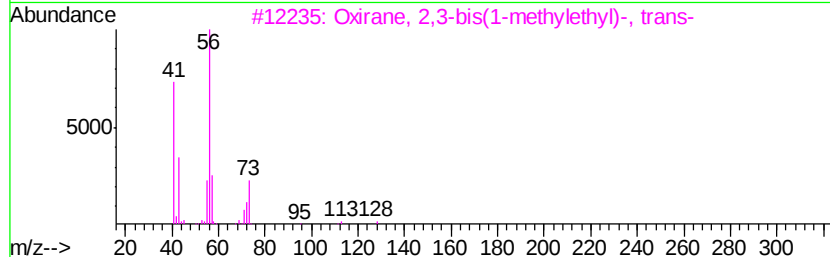
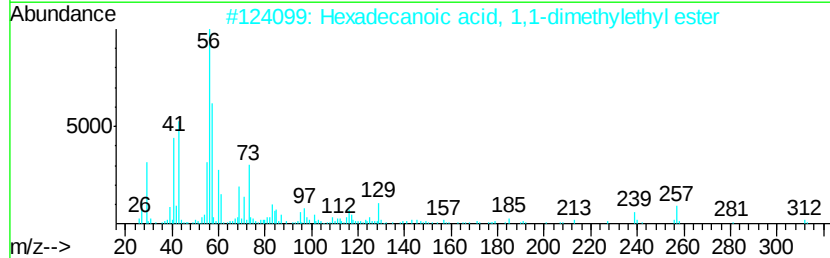
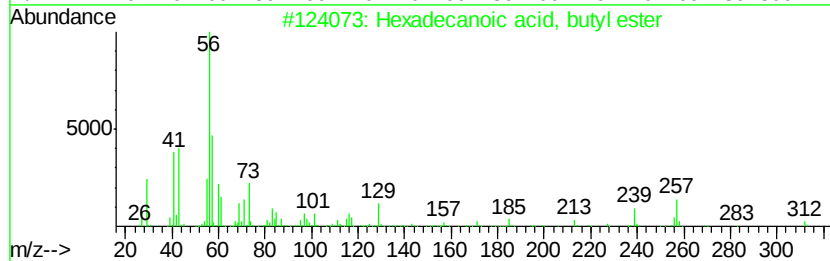
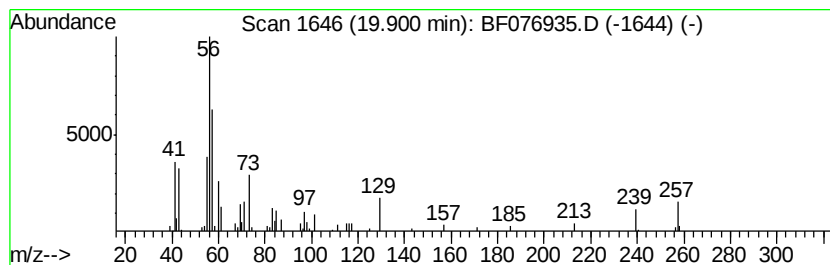
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecanoic acid, butyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.90	6.93 ng	107237	Chrysene-d12	21.27

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	87
3		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	37
4		1-Pentadecene, 2-methyl-	224	C16H32	029833-69-0	32
5		Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011915\
Data File : BF076935.D
Acq On : 19 Jan 2015 18:16
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Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

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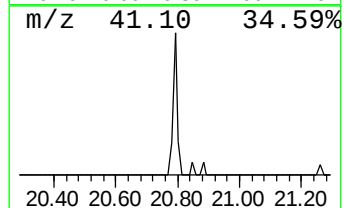
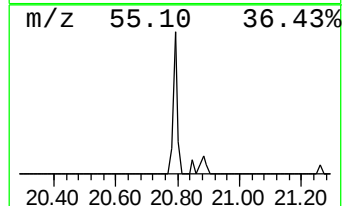
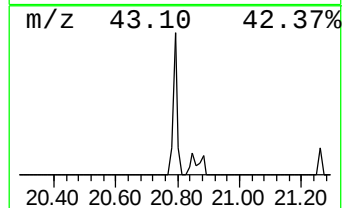
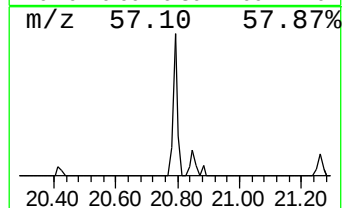
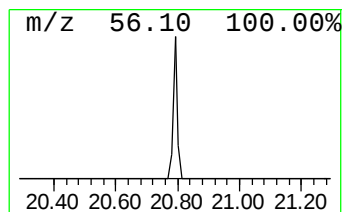
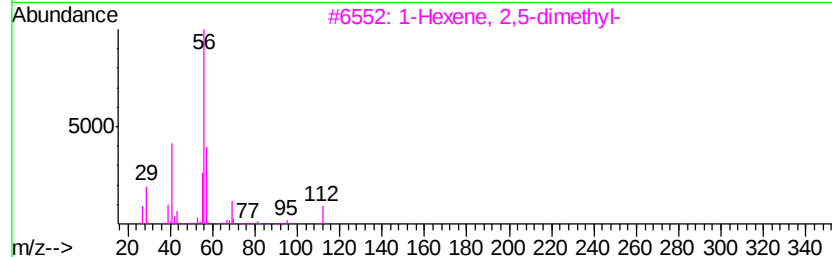
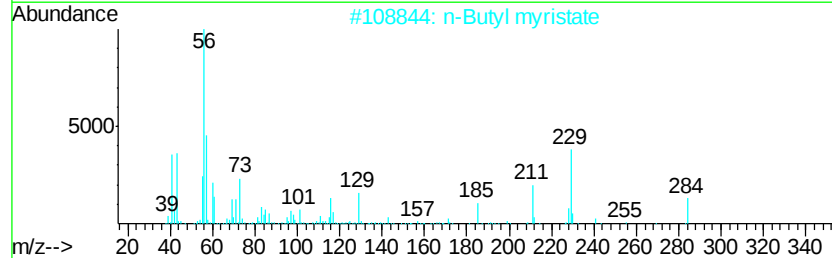
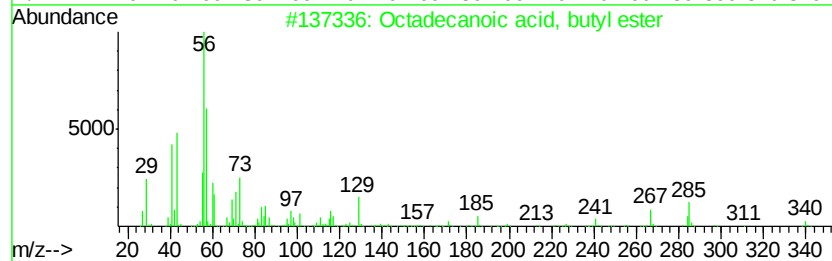
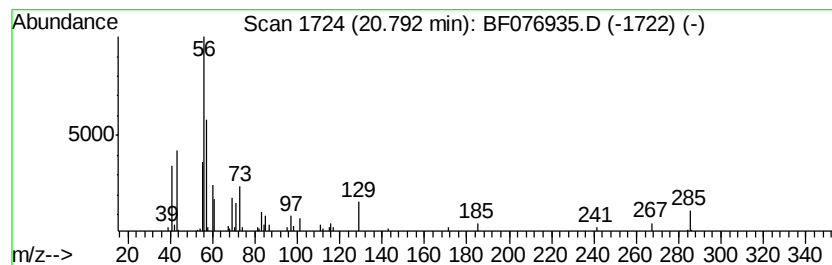
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octadecanoic acid, butyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	4.77 ng	73887	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	91
2		n-Butyl myristate	284	C18H36O2	000110-36-1	78
3		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	38
4		Pentadecane, 8-methylene-	224	C16H32	055668-09-2	37
5		1-Heptene, 2-methyl-	112	C8H16	015870-10-7	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011915\
Data File : BF076935.D
Acq On : 19 Jan 2015 18:16
Operator : IZ / TP
Sample : G1101-24
Misc : W
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FB-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.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
Oxirane, [(1-meth...	4.12	3.4	ng	33023	1	7.97	192393	20.0
unknown7.46	7.46	85.1	ng	819002	1	7.97	192393	20.0
Cycloheptanemethanol	11.40	2.9	ng	38250	2	10.88	261835	20.0
unknown12.14	12.14	3.7	ng	48435	2	10.88	261835	20.0
unknown13.96	13.96	18.7	ng	267261	3	15.01	285686	20.0
unknown14.76	14.76	4.4	ng	62563	3	15.01	285686	20.0
Benzene, (1,2-dim...	15.26	8.0	ng	114161	3	15.01	285686	20.0
Benzeneacetic aci...	15.36	4.8	ng	68022	3	15.01	285686	20.0
Bicyclo[4.2.0]oct...	16.31	3.6	ng	51457	3	15.01	285686	20.0
unknown16.47	16.47	32.5	ng	471534	4	17.77	289731	20.0
Hexadecanoic acid...	19.90	6.9	ng	107237	5	21.27	309690	20.0
Octadecanoic acid...	20.79	4.8	ng	73887	5	21.27	309690	20.0