

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
 Data File : BF079326.D
 Acq On : 19 May 2015 2:02
 Operator : TP/IZ
 Sample : G2271-15
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-13-D1

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	1.313	8	11	13	rVB	3384775	4033874	100.00%	11.779%
2	1.438	19	22	31	rVB	69006	170008	4.21%	0.496%
3	1.598	34	36	40	rVV	148070	194570	4.82%	0.568%
4	1.667	40	42	50	rVV	207052	353588	8.77%	1.032%
5	1.781	50	52	77	rVB	1647139	2847563	70.59%	8.315%
6	4.925	325	327	335	rVB	165396	249328	6.18%	0.728%
7	5.450	371	373	376	rBV	833321	1150753	28.53%	3.360%
8	6.742	483	486	490	rBV2	1282131	1948200	48.30%	5.689%
9	6.879	495	498	501	rBV	1441460	1324108	32.82%	3.866%
10	7.165	521	523	526	rBV	313985	295549	7.33%	0.863%
11	7.348	537	539	542	rBV	756004	971233	24.08%	2.836%
12	7.508	548	553	555	rBV2	109745	229313	5.68%	0.670%
13	7.690	567	569	572	rBV	40389	55849	1.38%	0.163%
14	7.862	581	584	586	rBV	892024	800720	19.85%	2.338%
15	8.342	625	626	629	rBV	57142	81848	2.03%	0.239%
16	8.445	633	635	638	rBV	2226291	2907725	72.08%	8.491%
17	8.742	659	661	662	rBV	485994	433650	10.75%	1.266%
18	8.765	662	663	665	rVV	329362	377110	9.35%	1.101%
19	9.245	703	705	709	rBV	94365	99466	2.47%	0.290%
20	9.485	723	726	728	rBV	707777	733576	18.19%	2.142%
21	9.634	730	739	740	rBV2	55306	144320	3.58%	0.421%
22	9.691	740	744	748	rVV	72024	193182	4.79%	0.564%
23	9.759	748	750	754	rVB2	32952	64595	1.60%	0.189%
24	9.862	754	759	763	rBV	33152	91643	2.27%	0.268%
25	9.942	763	766	768	rBV2	26572	44234	1.10%	0.129%
26	10.091	777	779	781	rVB	1866868	1586902	39.34%	4.634%
27	10.216	788	790	792	rVB	317244	387432	9.60%	1.131%
28	10.342	799	801	804	rVB	1114091	1065595	26.42%	3.112%
29	10.594	821	823	826	rBV	82378	92621	2.30%	0.270%
30	10.834	839	844	846	rVV2	46933	69959	1.73%	0.204%
31	10.914	849	851	856	rVB2	485085	679528	16.85%	1.984%
32	11.497	897	902	904	rBV	24112	43750	1.08%	0.128%
33	11.565	906	908	913	rVV3	30017	49701	1.23%	0.145%
34	11.897	934	937	940	rBV	1088063	1094650	27.14%	3.196%

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Sampling : 1 Min Area: 3 % of largest Peak
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35	12.091	952	954	959	rBV	39903	55766	1.38%	0.163%
36	12.754	1010	1012	1017	rBV2	493588	602876	14.95%	1.760%
37	12.845	1017	1020	1023	rBV	44327	60965	1.51%	0.178%
38	13.474	1072	1075	1077	rBV2	28680	55133	1.37%	0.161%
39	13.520	1077	1079	1081	rVV2	34868	47584	1.18%	0.139%
40	13.794	1101	1103	1106	rBV	259292	349649	8.67%	1.021%
41	14.091	1125	1129	1130	rBV2	130361	194924	4.83%	0.569%
42	14.125	1130	1132	1134	rVB	999449	1156657	28.67%	3.377%
43	14.263	1142	1144	1146	rBV	205002	206410	5.12%	0.603%
44	14.445	1157	1160	1162	rBV	612184	649295	16.10%	1.896%
45	14.537	1166	1168	1170	rVV	179679	229998	5.70%	0.672%
46	14.708	1181	1183	1184	rBV	82214	125620	3.11%	0.367%
47	14.754	1184	1187	1190	rVB	1855292	1908399	47.31%	5.573%
48	14.971	1203	1206	1209	rVV3	83661	184972	4.59%	0.540%
49	15.017	1209	1210	1215	rVB2	90301	116077	2.88%	0.339%
50	15.234	1228	1229	1231	rBV	43922	46026	1.14%	0.134%
51	15.280	1231	1233	1238	rVV3	23242	66099	1.64%	0.193%
52	15.920	1286	1289	1293	rBV2	210686	328160	8.14%	0.958%
53	16.046	1296	1300	1301	rVV2	471612	751383	18.63%	2.194%
54	16.091	1301	1304	1306	rVB	580630	838152	20.78%	2.447%
55	16.331	1323	1325	1330	rBV6	22204	41665	1.03%	0.122%
56	16.686	1354	1356	1357	rBV2	26456	41515	1.03%	0.121%
57	16.720	1357	1359	1363	rVV3	37724	108345	2.69%	0.316%
58	17.303	1408	1410	1416	rBV	122195	270196	6.70%	0.789%
59	17.611	1435	1437	1440	rBV	89625	124718	3.09%	0.364%
60	17.669	1440	1442	1445	rVV	119205	176014	4.36%	0.514%
61	17.737	1445	1448	1455	rVB	469906	643597	15.95%	1.879%

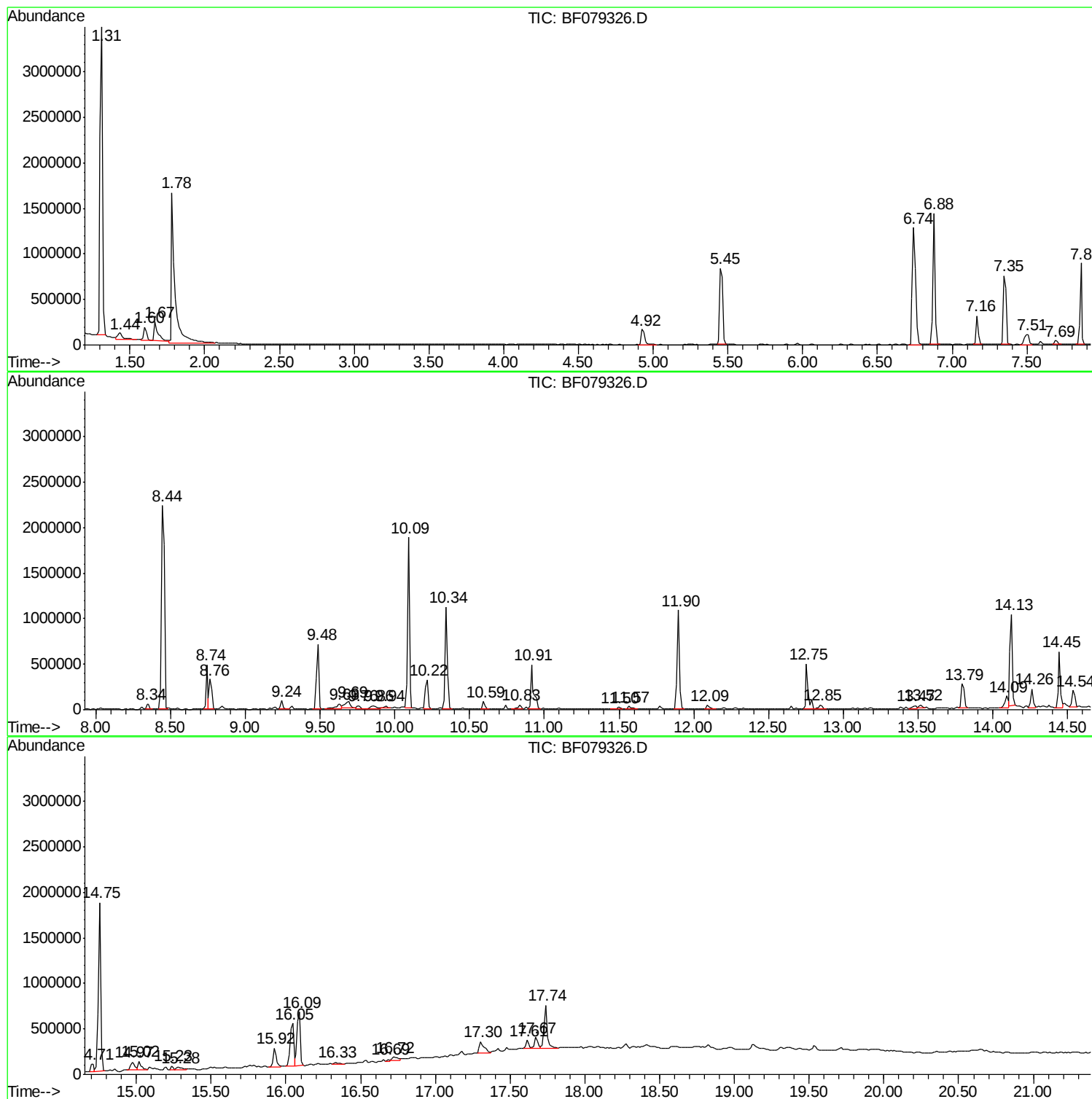
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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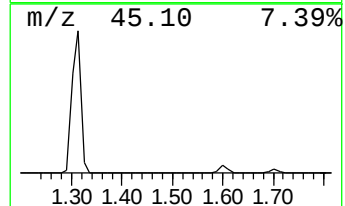
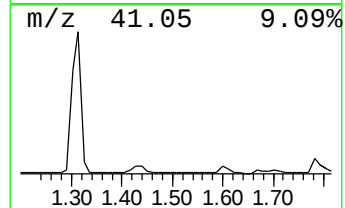
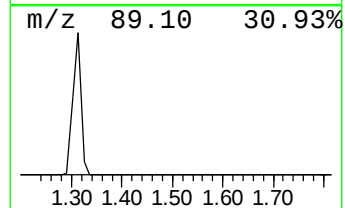
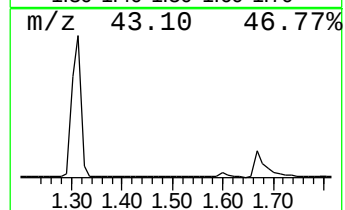
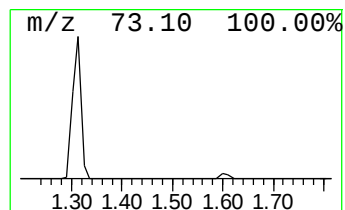
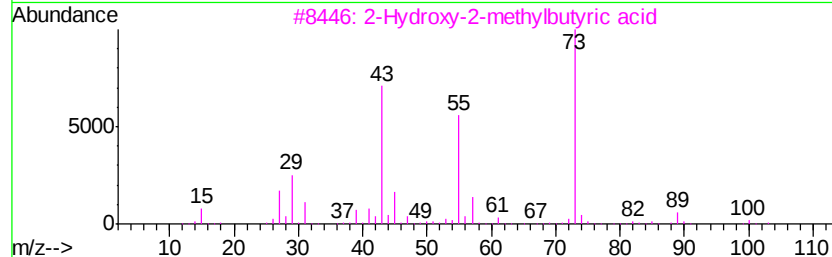
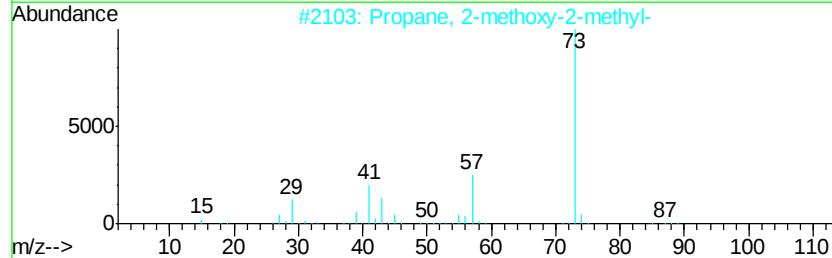
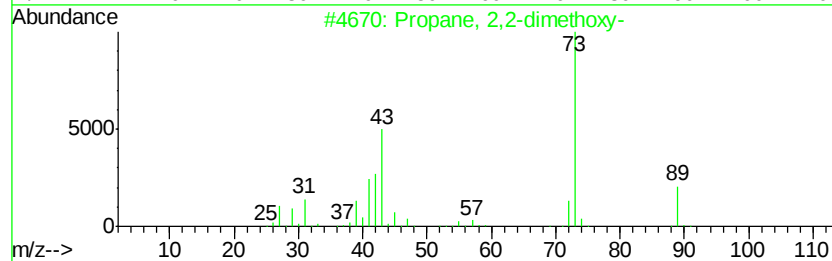
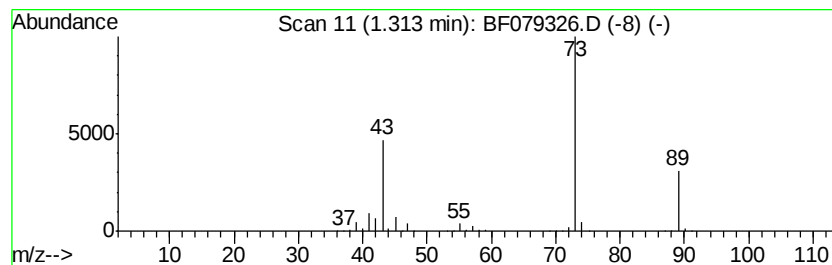
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.31 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.31	272.98 ng	4033870	1,4-Dichlorobenzene-d4	7.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9
5		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	7



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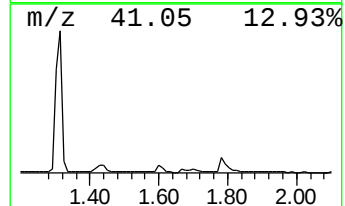
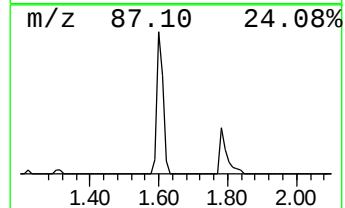
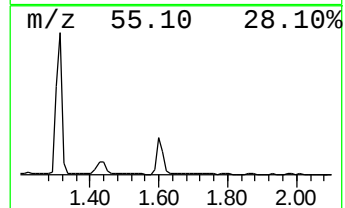
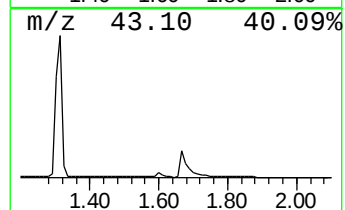
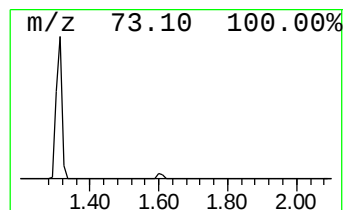
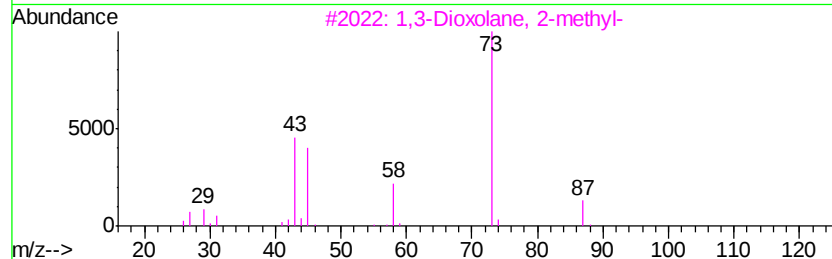
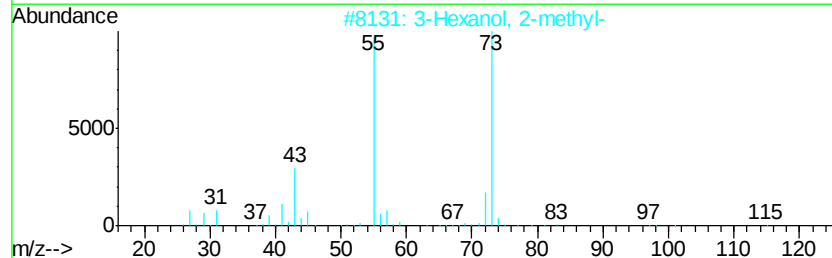
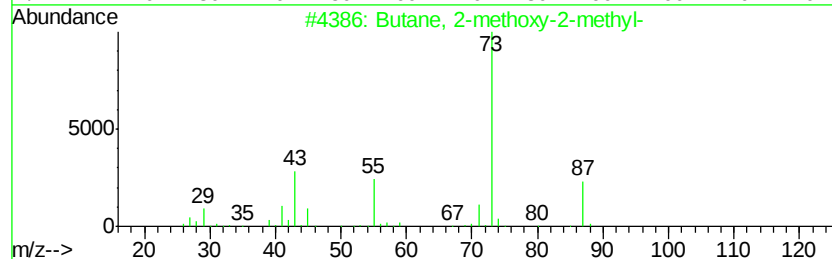
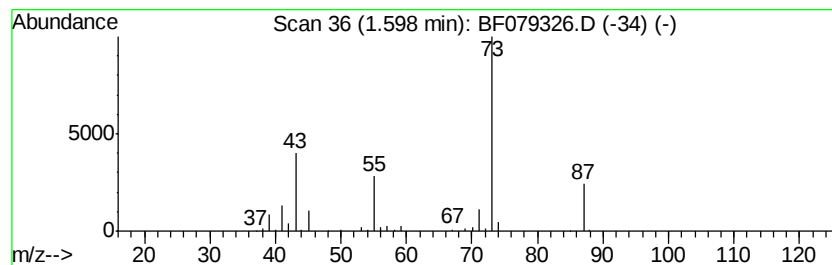
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.60	13.17 ng	194570	1,4-Dichlorobenzene-d4	7.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	25
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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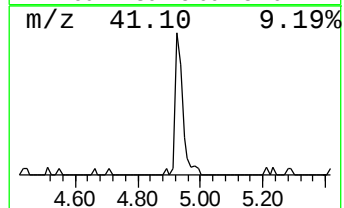
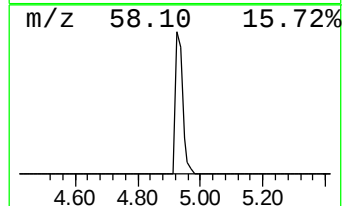
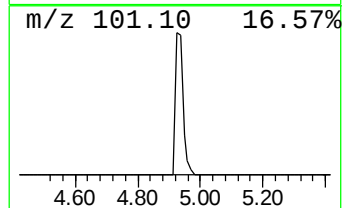
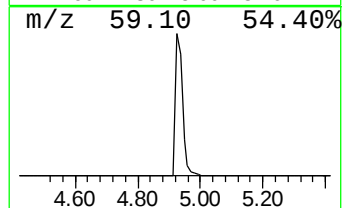
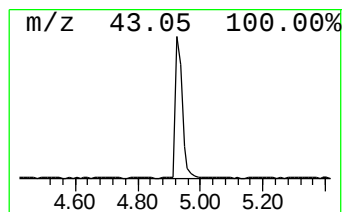
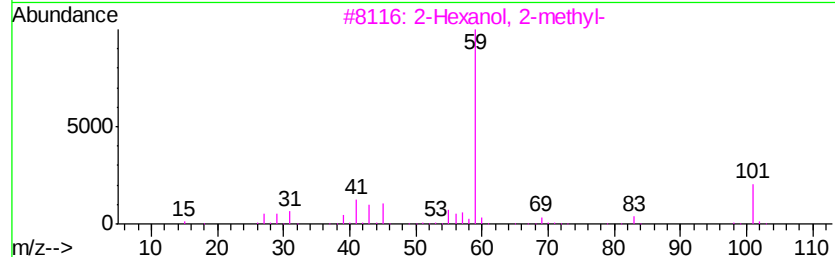
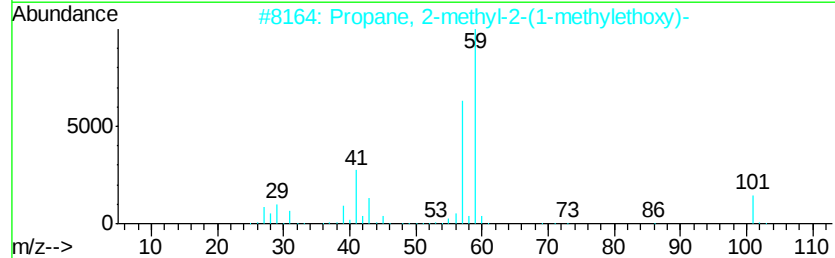
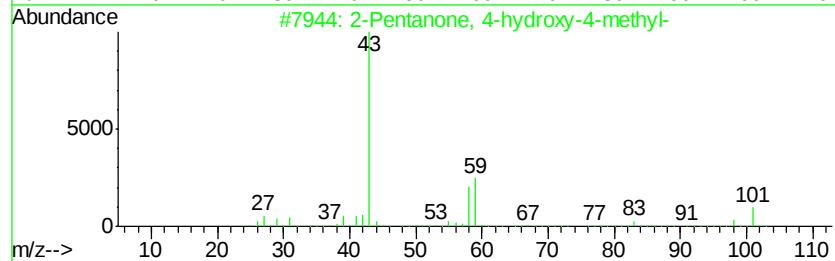
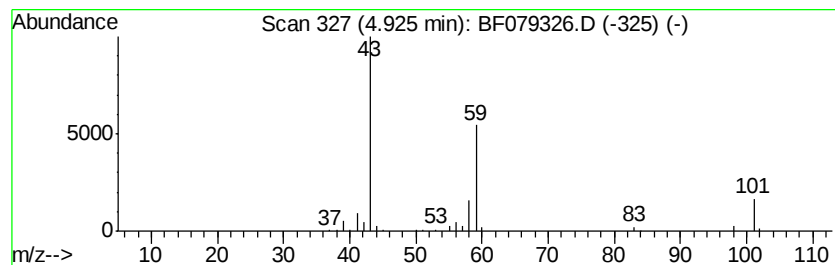
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	16.87 ng	249328	1,4-Dichlorobenzene-d4	7.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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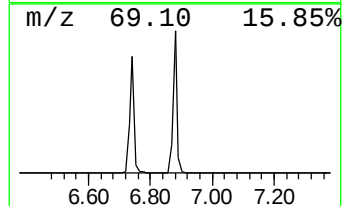
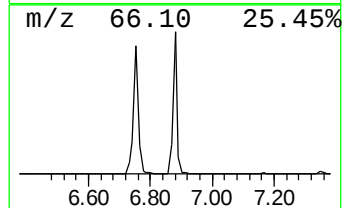
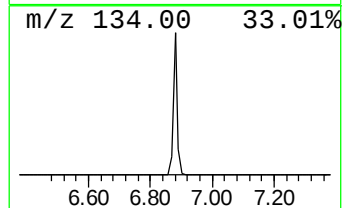
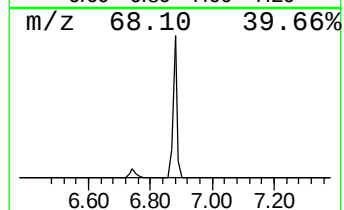
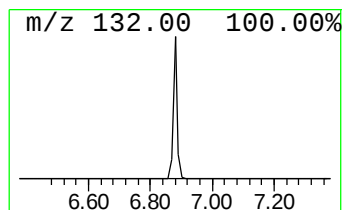
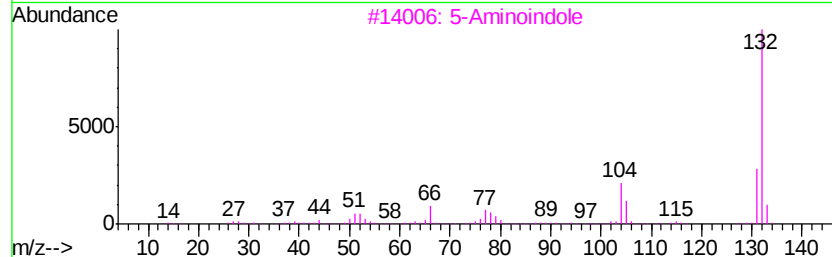
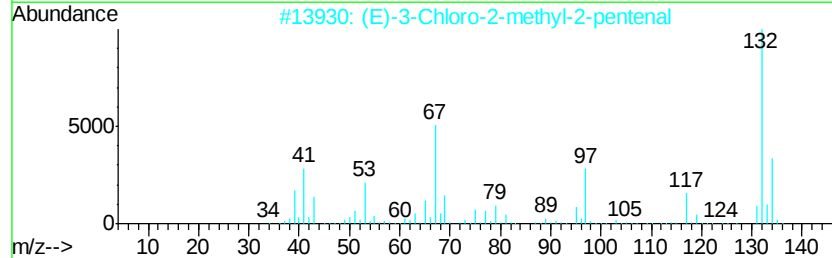
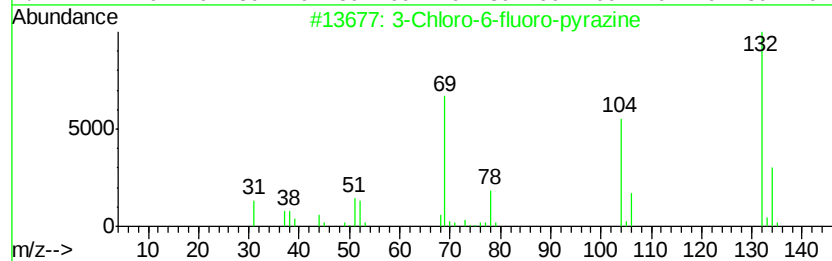
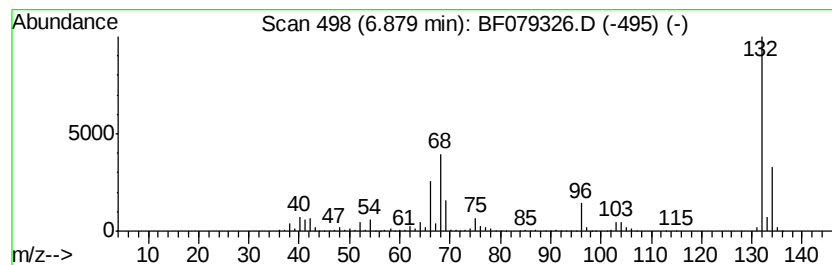
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown6.88 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	89.60 ng	1324110	1,4-Dichlorobenzene-d4	7.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	17
3	5-Aminoindole			132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
5	Amphetaminil			250	C17H18N2	017590-01-1	10



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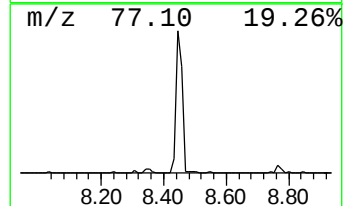
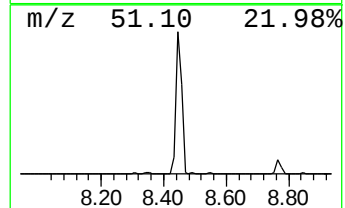
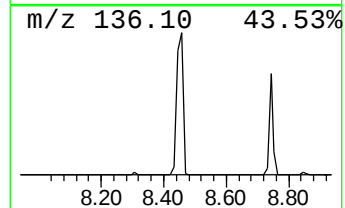
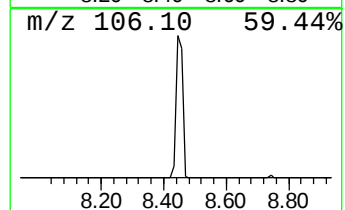
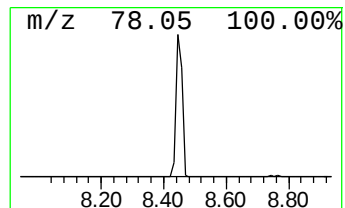
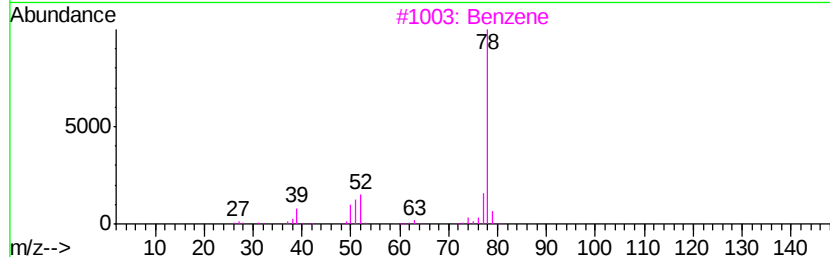
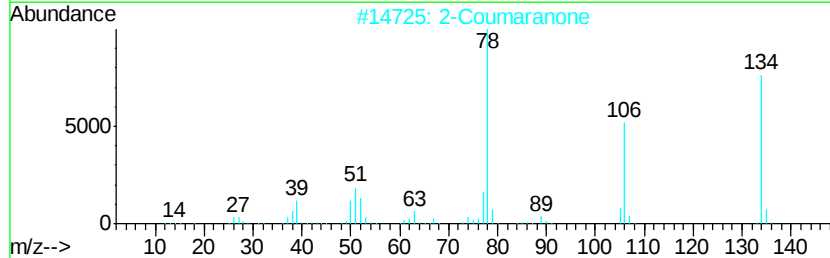
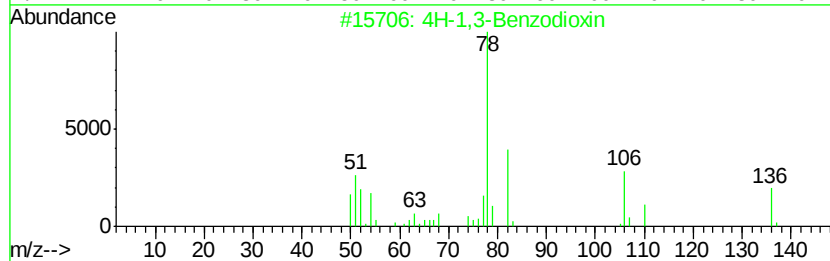
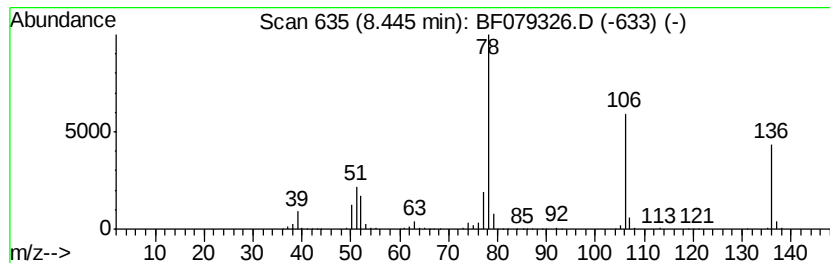
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-1,3-Benzodioxin Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.44	134.11 ng	2907730	Naphthalene-d8	8.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-1,3-Benzodioxin	136	C8H8O2	000254-27-3	86
2		2-Coumaranone	134	C8H6O2	000553-86-6	72
3		Benzene	78	C6H6	000071-43-2	68
4		2-Nitrobenzyl bromide	215	C7H6BrNO2	003958-60-9	56
5		Ethanone, 1-(3-pyridinyl)-	121	C7H7NO	000350-03-8	56



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
Operator : TP/IZ
Sample : G2271-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleID :
WC-13-D1

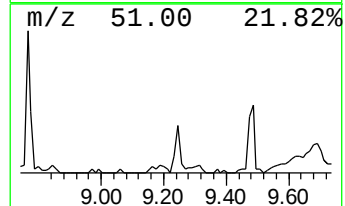
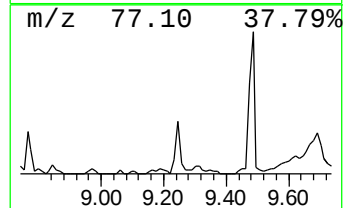
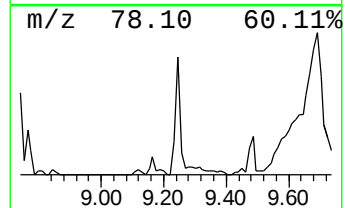
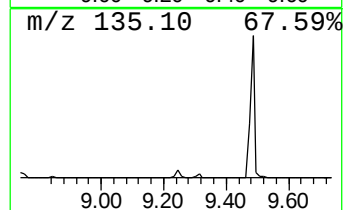
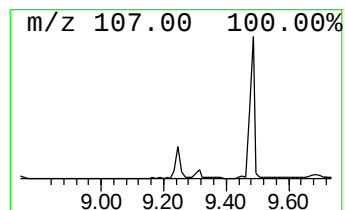
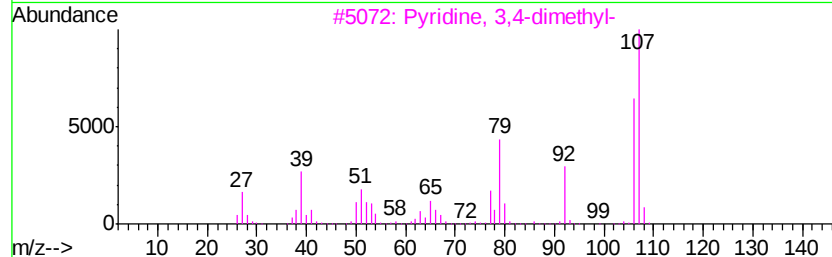
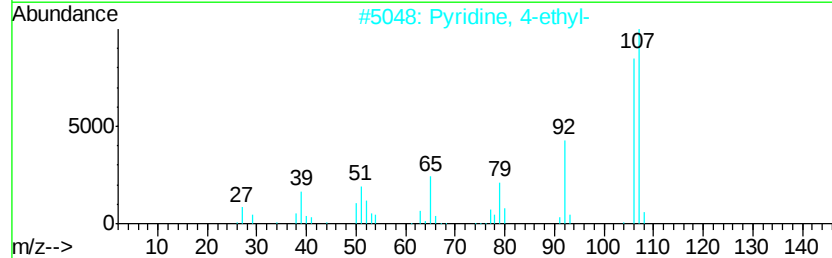
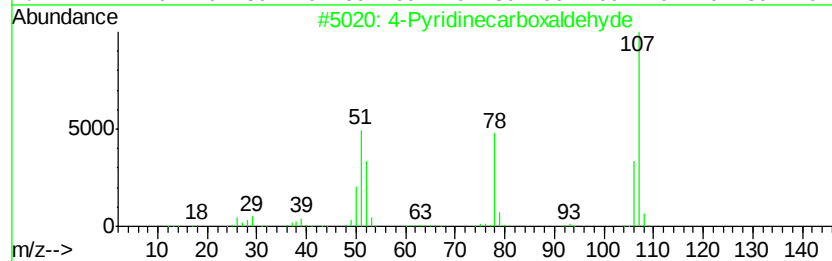
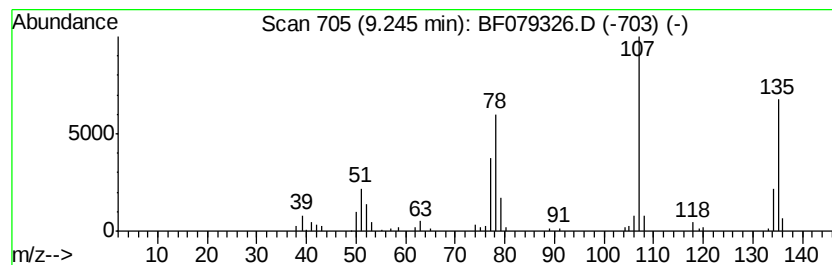
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.24 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.24	4.59 ng	99466	Naphthalene-d8	8.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Pyridinecarboxaldehyde	107	C6H5NO	000872-85-5	47
2			Pyridine, 4-ethyl-	107	C7H9N	000536-75-4	35
3			Pyridine, 3,4-dimethyl-	107	C7H9N	000583-58-4	25
4			2-Butenedinitrile, (E)-	78	C4H2N2	000764-42-1	18
5			Ethanedione, (4-methoxyphenyl)ph...	240	C15H12O3	022711-21-3	12



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
Operator : TP/IZ
Sample : G2271-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-13-D1

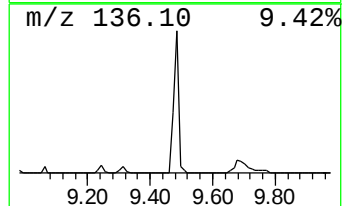
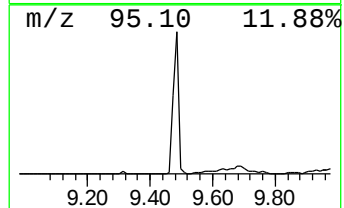
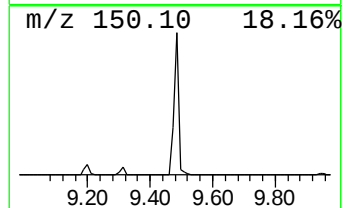
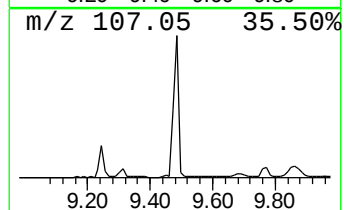
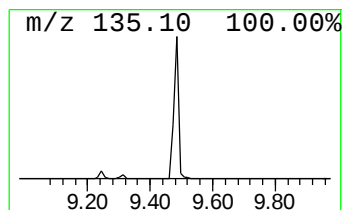
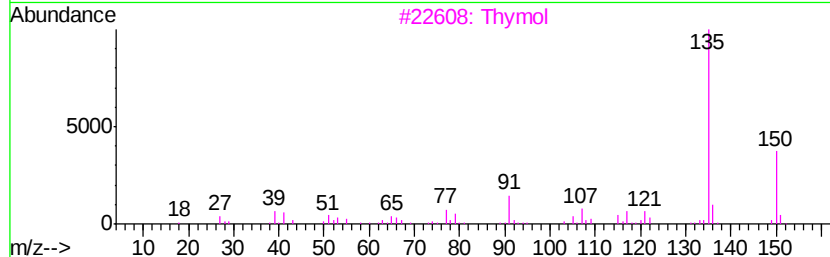
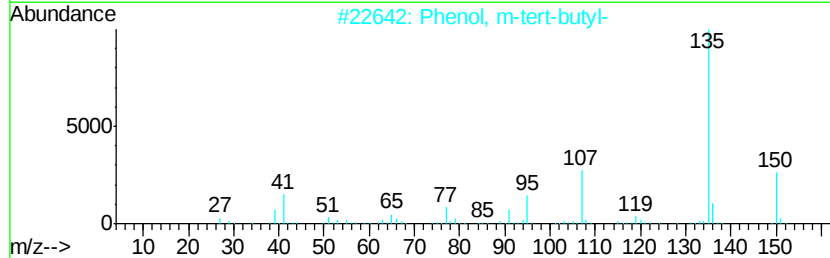
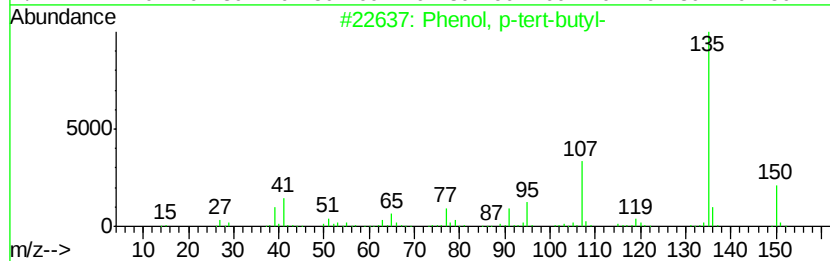
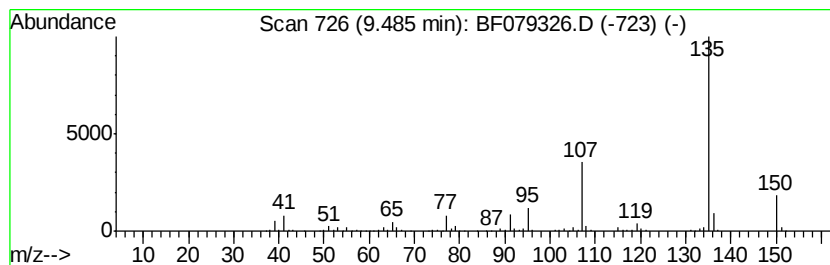
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenol, p-tert-butyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	33.83 ng	733576	Napthalene-d8	8.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
2		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	95
3		Thymol	150	C10H14O	000089-83-8	72
4		Phenol, 2,3,5,6-tetramethyl-	150	C10H14O	000527-35-5	59
5		Ethanone, 1-(3-methoxyphenyl)-	150	C9H10O2	000586-37-8	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
Operator : TP/IZ
Sample : G2271-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

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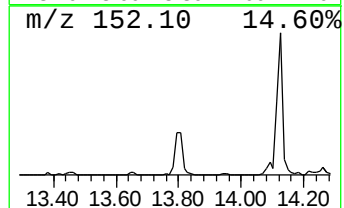
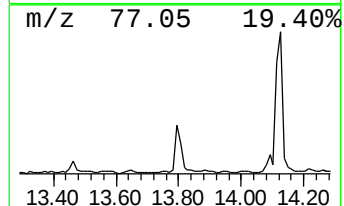
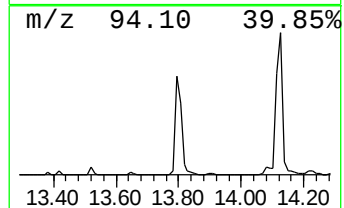
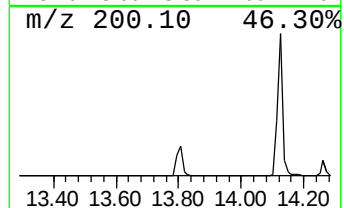
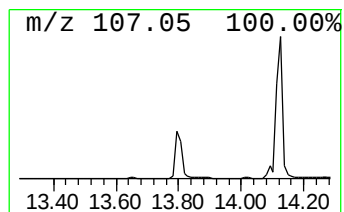
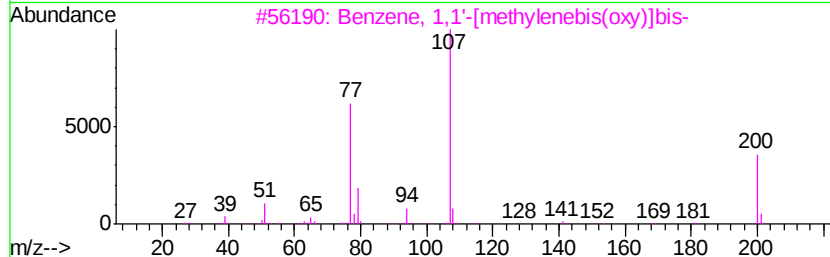
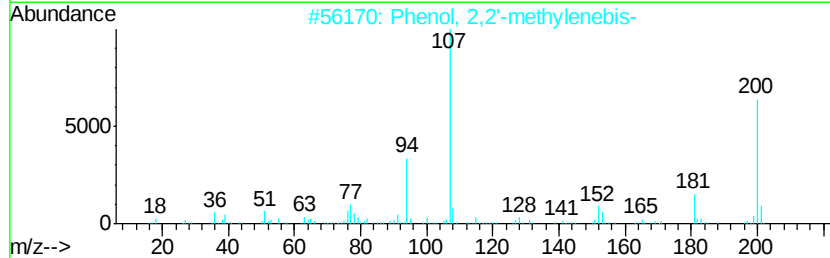
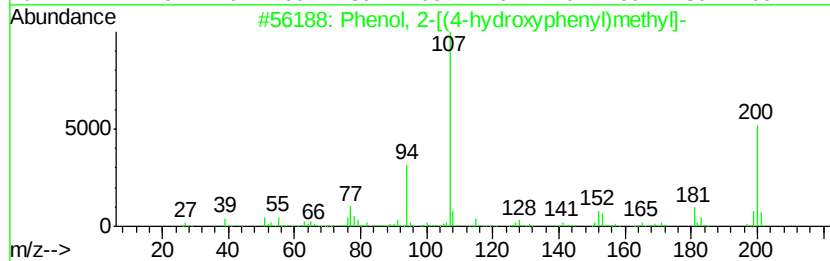
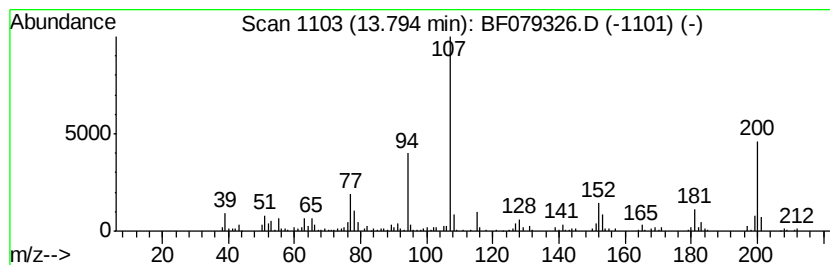
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Phenol, 2-[(4-hydroxyphenyl)... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	11.60 ng	349649	Phenanthrene-d10	12.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2-[(4-hydroxyphenyl)meth...	200	C13H12O2	002467-03-0	97
2		Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	90
3		Benzene, 1,1'-[methylenebis(oxy)]...	200	C13H12O2	004442-41-5	47
4		4-(2-Methoxyethyl)phenol	152	C9H12O2	056718-71-9	43
5		Benzeneacetic acid, 4-hydroxy-	152	C8H8O3	000156-38-7	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
Operator : TP/IZ
Sample : G2271-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WC-13-D1

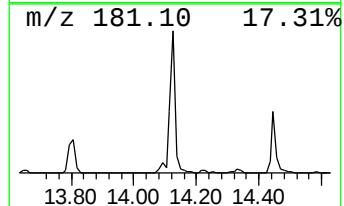
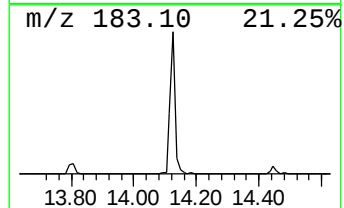
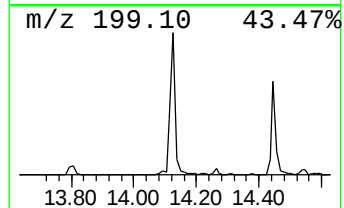
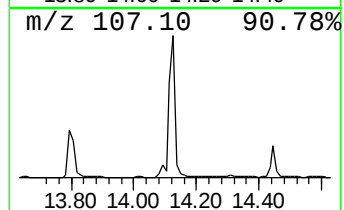
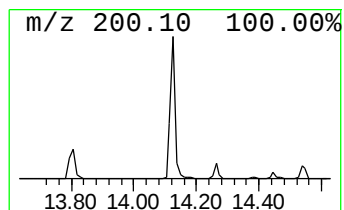
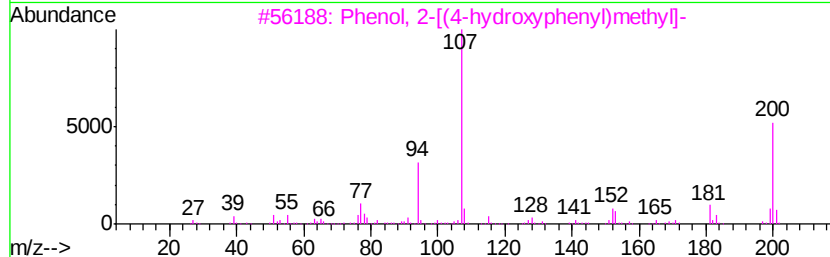
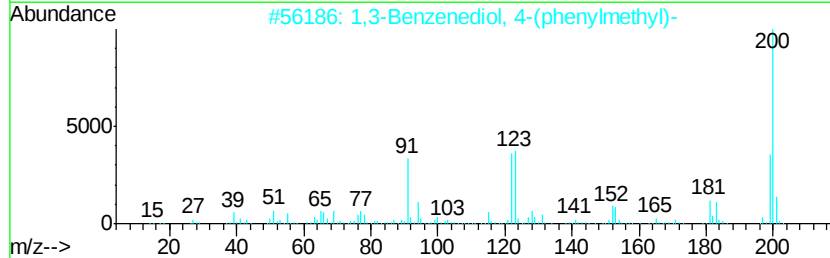
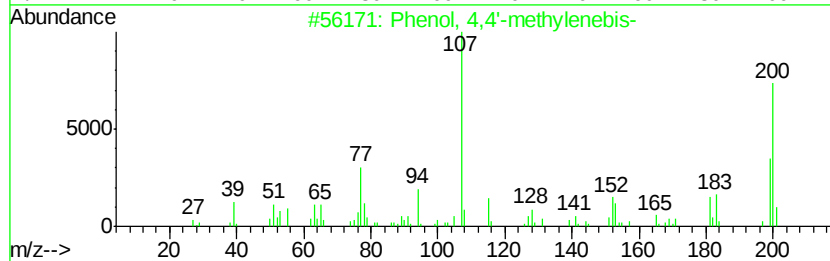
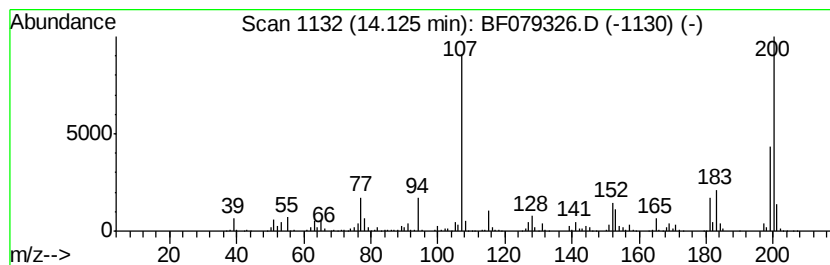
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Phenol, 4,4'-methylenebis- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	38.37 ng	1156660	Phenanthrene-d10	12.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-methylenebis-	200	C13H12O2	000620-92-8	98
2		1,3-Benzenediol, 4-(phenylmethyl)-	200	C13H12O2	002284-30-2	83
3		Phenol, 2-[(4-hydroxyphenyl)meth]...	200	C13H12O2	002467-03-0	81
4		Phenol, 2,2'-methylenebis-	200	C13H12O2	002467-02-9	81
5		Thiazolo[5,4-f]quinoline, 7-methyl-	200	C11H8N2S	003119-56-0	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
Operator : TP/IZ
Sample : G2271-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
WC-13-D1

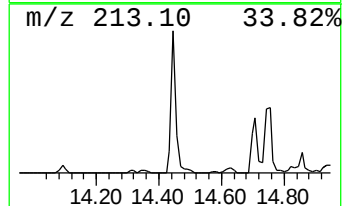
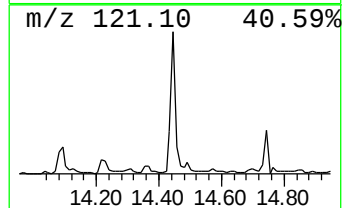
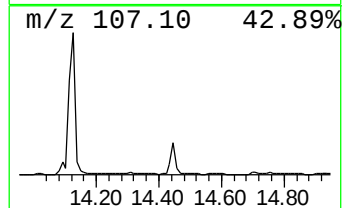
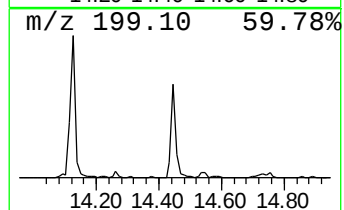
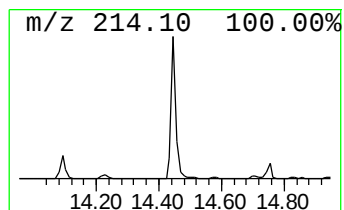
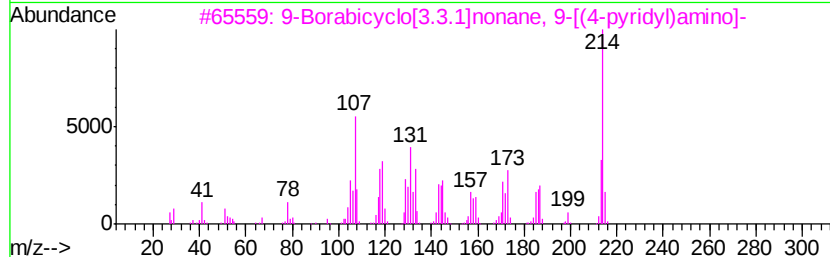
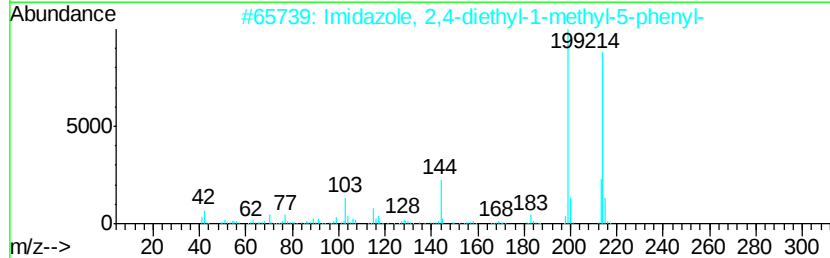
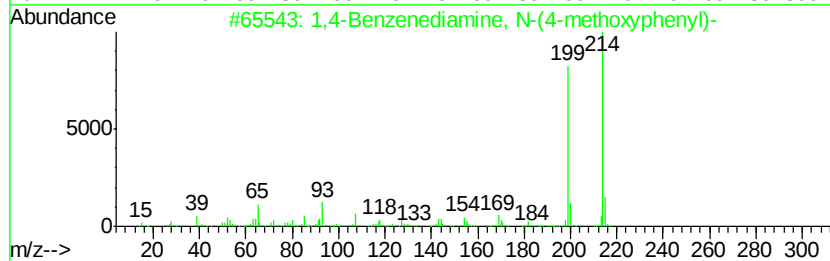
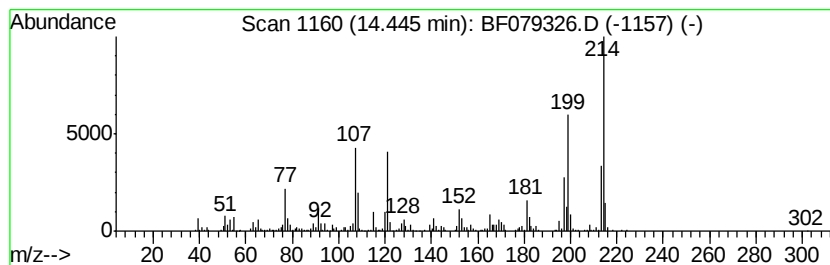
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown14.45 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	17.28 ng	649295	Chrysene-d12	16.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Benzenediamine, N-(4-methoxy...	214	C13H14N2O	000101-64-4	49
2		Imidazole, 2,4-diethyl-1-methyl-...	214	C14H18N2	062576-17-4	47
3		9-Borabicyclo[3.3.1]nonane, 9-[(...	214	C13H19BN2	1000158-11-2	43
4		4,4'-Ethylidenediphenol	214	C14H14O2	002081-08-5	43
5		3,4,5,6-Tetramethoxyphenol	214	C10H14O5	030225-99-1	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
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Misc :
ALS Vial : 14 Sample Multiplier: 1

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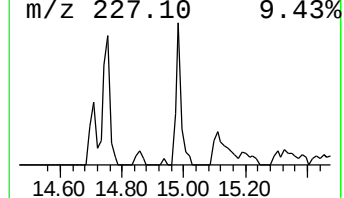
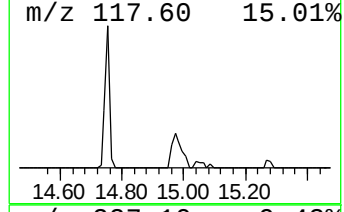
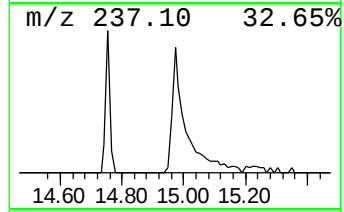
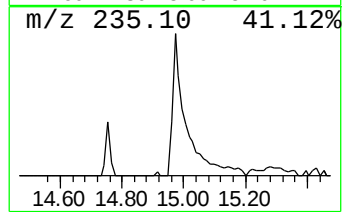
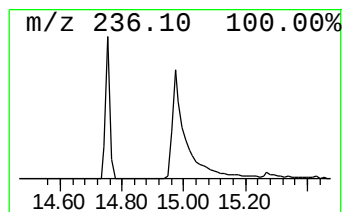
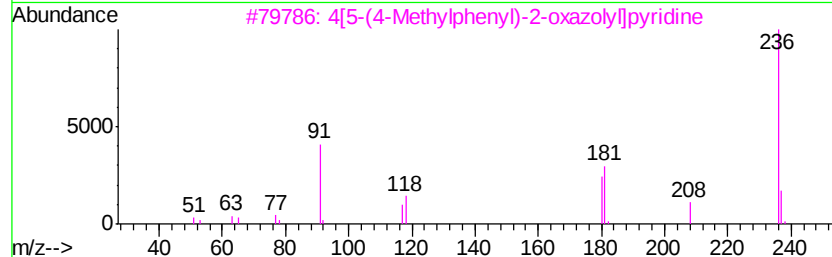
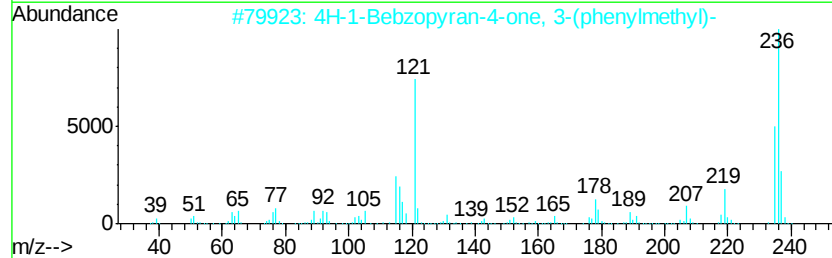
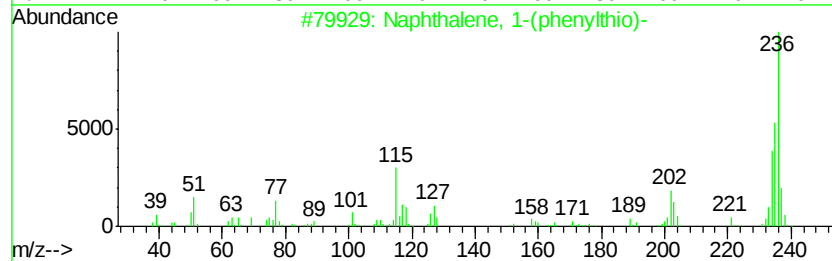
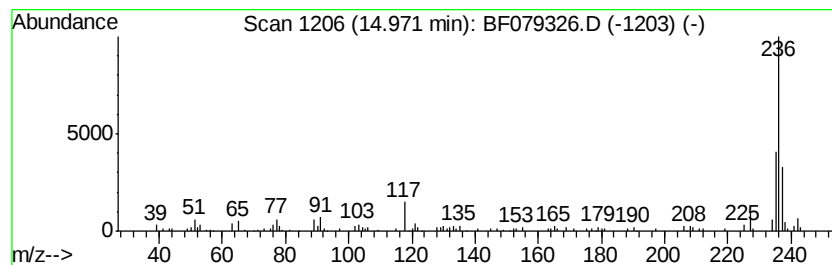
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Naphthalene, 1-(phenylthio)- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	4.92 ng	184972	Chrysene-d12	16.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-(phenylthio)-	236	C16H12S	007570-98-1	64
2		4H-1-Benzopyran-4-one, 3-(phenyl...)	236	C16H12O2	030779-85-2	64
3		4[5-(4-Methylphenyl)-2-oxazolyl]...	236	C15H12N2O	097813-94-0	46
4		Pyrido[2,3-d]pyrimidin-5(8H)-one...	237	C14H11N3O	161466-00-8	46
5		1H-Pyrazole, 4,5-dihydro-1-(4-me...	236	C16H16N2	000959-08-0	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
Data File : BF079326.D
Acq On : 19 May 2015 2:02
Operator : TP/IZ
Sample : G2271-15
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-13-D1

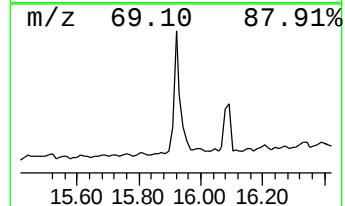
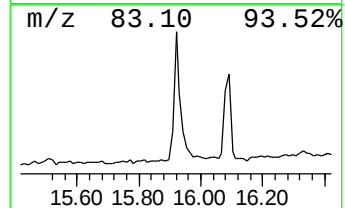
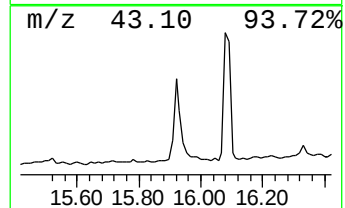
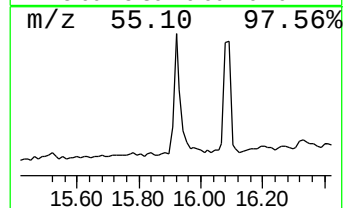
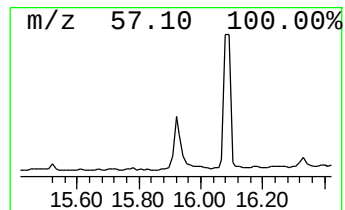
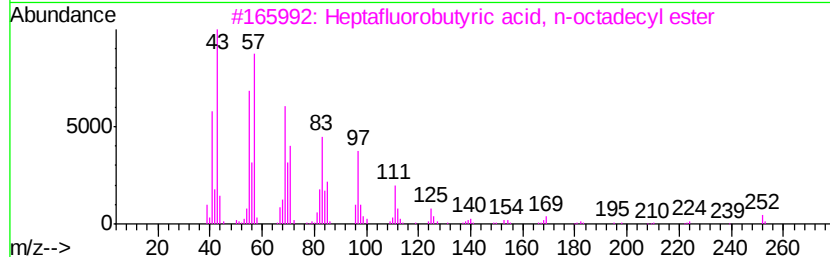
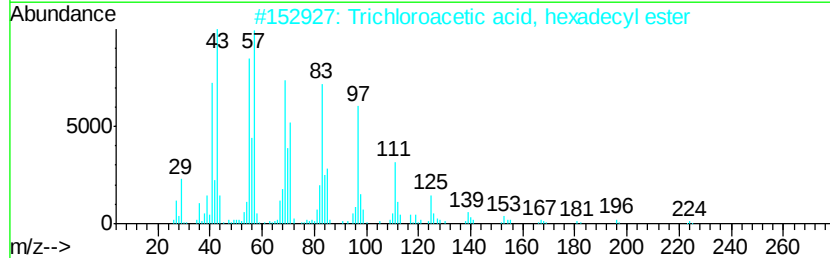
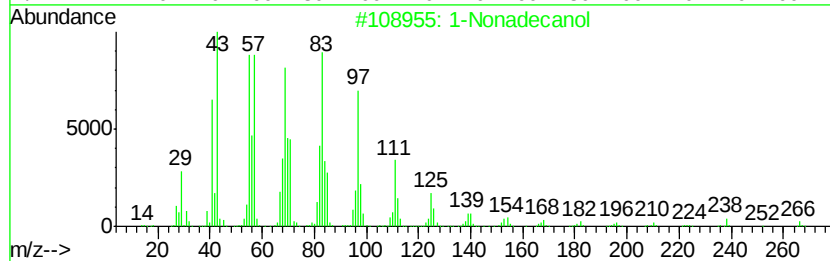
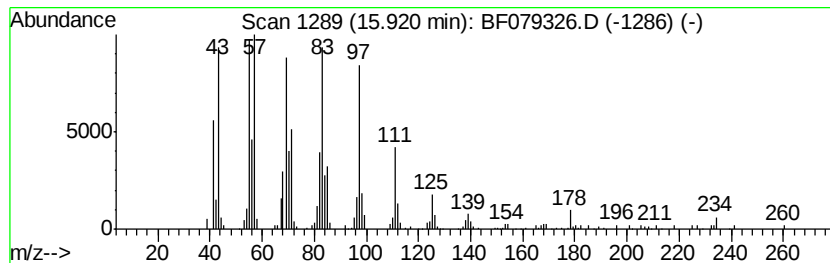
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 1-Nonadecanol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	8.73 ng	328160	Chrysene-d12	16.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecanol	284	C19H40O	001454-84-8	95
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
3		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
4		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
5		1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051815\
 Data File : BF079326.D
 Acq On : 19 May 2015 2:02
 Operator : TP/IZ
 Sample : G2271-15
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-13-D1

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
unknown1.31	1.31	273.0	ng	4033870	1	7.16	295549	20.0
Butane, 2-methoxy...	1.60	13.2	ng	194570	1	7.16	295549	20.0
2-Pentanone, 4-hy...	4.92	16.9	ng	249328	1	7.16	295549	20.0
unknown6.88	6.88	89.6	ng	1324110	1	7.16	295549	20.0
4H-1,3-Benzodioxin	8.44	134.1	ng	2907730	2	8.74	433650	20.0
unknown9.24	9.24	4.6	ng	99466	2	8.74	433650	20.0
Phenol, p-tert-bu...	9.48	33.8	ng	733576	2	8.74	433650	20.0
Phenol, 2-[(4-hyd...	13.79	11.6	ng	349649	4	12.75	602876	20.0
Phenol, 4,4'-meth...	14.13	38.4	ng	1156660	4	12.75	602876	20.0
unknown14.45	14.45	17.3	ng	649295	5	16.05	751383	20.0
Naphthalene, 1-(p...	14.97	4.9	ng	184972	5	16.05	751383	20.0
1-Nonadecanol	15.92	8.7	ng	328160	5	16.05	751383	20.0