LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080156.D
Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

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-BF061715.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	5.533	211	214	216	rBV	310348	345695	21.62%	2.538%
2	5.921	246	248	251	rBV	922475	1020023	63.78%	7.488%
3	6.161	267	269	276	rVB2	12454	25459	1.59%	0.187%
4	6.321	281	283	285	rBV	30731	27653	1.73%	0.203%
5	6.550	300	303	307	rBV	13506	21567	1.35%	0.158%
6	6.916	333	335	338	rBV	1179373	1017361	63.61%	7.468%
7	7.076	347	349	352	rVV	853443	1060186	66.29%	7.783%
8	7.144	352	355	357	rVB2	14321	16595	1.04%	0.122%
9	7.316	368	370	372	rBV	249108	196801	12.31%	1.445%
10	7.476	381	384	386	rBV	1056820	864462	54.05%	6.346%
11	7.876	416	419	421	rBV	456951	624655	39.06%	4.586%
12	8.367	459	462	464	rVB	18381	19613	1.23%	0.144%
13	8.607	481	483	484	rBV	326022	267720	16.74%	1.965%
14	8.630	484	485	487	rVB	30327	21090	1.32%	0.155%
15	8.847	503	504	506	rBV	24011	18318	1.15%	0.134%
16	8.939	510	512	513	rBV	117412	109720	6.86%	0.805%
17	8.962	513	514	516	rVV	26897	19434	1.22%	0.143%
18	9.064	517	523	525	rBV	67728	107364	6.71%	0.788%
19	9.270	537	541	544	rBV	35992	48908	3.06%	0.359%
20	9.316	544	545	547	rVB	24785	29474	1.84%	0.216%
21	9.350	547	548	550	rBV	20286	27660	1.73%	0.203%
22	9.442	554	556	559	rVV2	34308	51294	3.21%	0.377%
23	9.556	563	566	571	rVB2	11687	28185	1.76%	0.207%
24	9.682	575	577	579	rVB	1611732	1266754	79.21%	9.299%
25	10.070	609	611	613	rBV	132332	105432	6.59%	0.774%
26	10.265	624	628	633	rVB2	7158	16190	1.01%	0.119%
27	10.379	635	638	640	rBV	370781	333727	20.87%	2.450%
28	11.168	705	707	709	rBV	849473	838767	52.45%	6.157%
29	11.259	714	715	721	rVB	9374	16011	1.00%	0.118%
30	11.396	724	727	729	rVB	21194	21312	1.33%	0.156%
31	11.876	767	769	773	rBV	323300	368992	23.07%	2.709%
32	12.082	785	787	791	rBV	16453	25487	1.59%	0.187%
33	12.596	828	832	835	rVB4	18069	23375	1.46%	0.172%
34	12.962	863	864	867	rBV2	7770	18467	1.15%	0.136%

LSC Area Percent Report

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Data File : BF080156.D
Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

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-BF061715.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.076	873	874	877	rBV3	18309	16261	1.02%	0.119%
36	13.145	878	880	882	rBV	131190	111411	6.97%	0.818%
37	13.316	893	895	897	rBV2	16396	22833	1.43%	0.168%
38	13.396	898	902	905	rVB	144731	159158	9.95%	1.168%
39	13.545	912	915	917	rBV	1215630	1599313	100.00%	11.740%
40	13.751	929	933	937	rBV2	17603	37919	2.37%	0.278%
41	13.831	938	940	942	rVB	15750	17406	1.09%	0.128%
42	13.876	942	944	945	rBV2	13237	18992	1.19%	0.139%
43	14.036	955	958	960	rBV2	15691	22845	1.43%	0.168%
44	14.482	994	997	998	rBV	12451	18838	1.18%	0.138%
45	14.562	1001	1004	1009	rVV4	157427	246658	15.42%	1.811%
46	14.734	1017	1019	1020	rBV	19057	19162	1.20%	0.141%
47	14.779	1020	1023	1025	rVV2	415849	598852	37.44%	4.396%
48	14.814	1025	1026	1029	rVB	73660	60971	3.81%	0.448%
49	14.905	1032	1034	1036	rBV3	19451	21682	1.36%	0.159%
50	14.962	1037	1039	1041	rBV2	12820	18074	1.13%	0.133%
51	15.054	1044	1047	1048	rBV2	11448	17425	1.09%	0.128%
52	15.259	1063	1065	1067	rVB	17493	21769	1.36%	0.160%
53	15.362	1071	1074	1078	rBV2	125522	192438	12.03%	1.413%
54	15.568	1090	1092	1094	rVB2	16085	22427	1.40%	0.165%
55	16.082	1134	1137	1143	rBV	102395	257030	16.07%	1.887%
56	16.174	1143	1145	1147	rBV	15527	27521	1.72%	0.202%
57	16.219	1147	1149	1151	rVB	27575	40520	2.53%	0.297%
58	16.459	1167	1170	1173	rVB	71696	103703	6.48%	0.761%
59	16.528	1174	1176	1181	rBV	83864	147824	9.24%	1.085%
60	16.619	1181	1184	1186	rBV	257400	438176	27.40%	3.217%
61	17.271	1239	1241	1247	rVB5	21360	61225	3.83%	0.449%
62	18.448	1340	1344	1350	rVB2	58654	145298	9.09%	1.067%
63	18.997	1388	1392	1401	rVB	48327	152722	9.55%	1.121%

Sum of corrected areas: 13622204

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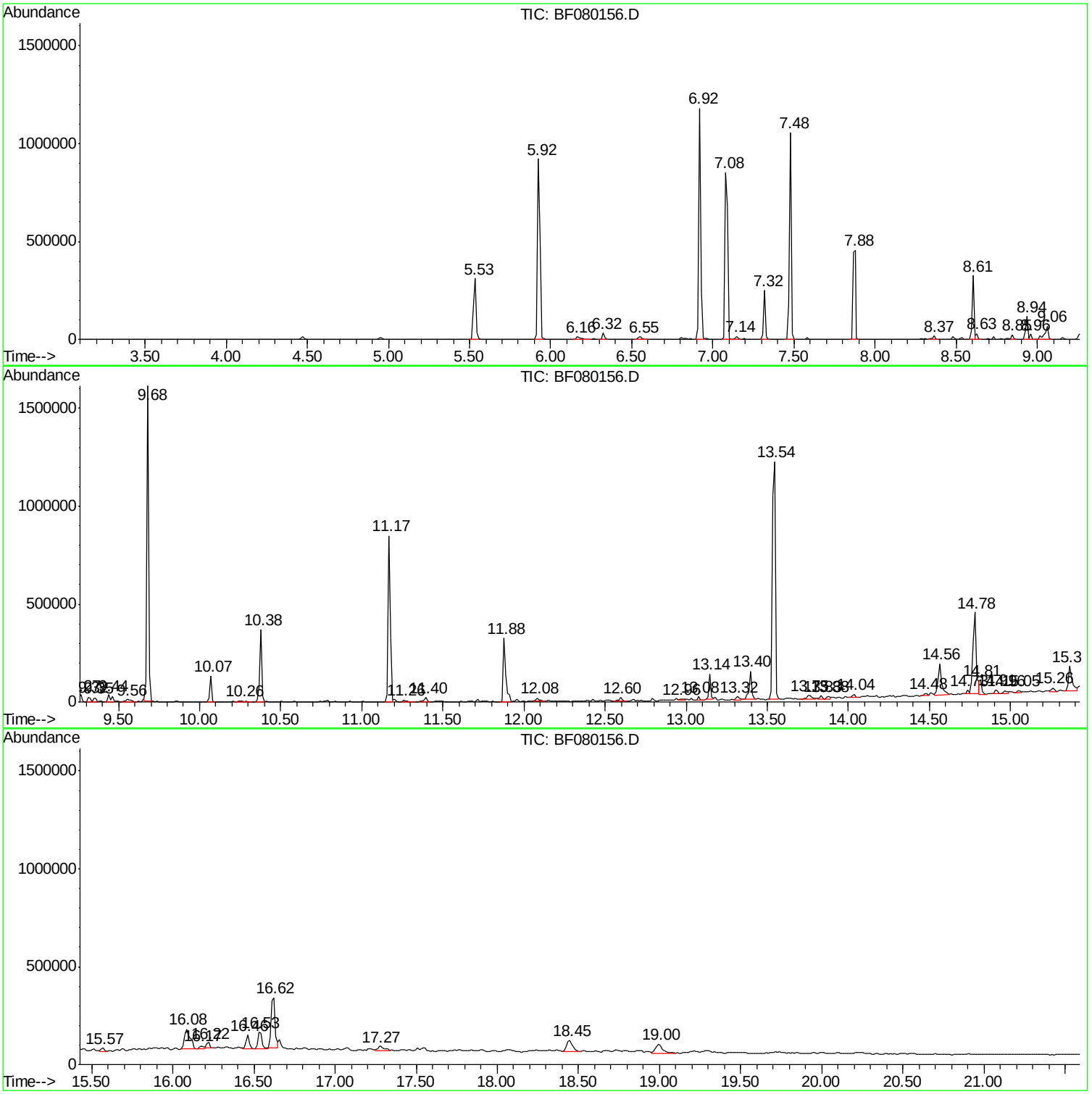
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

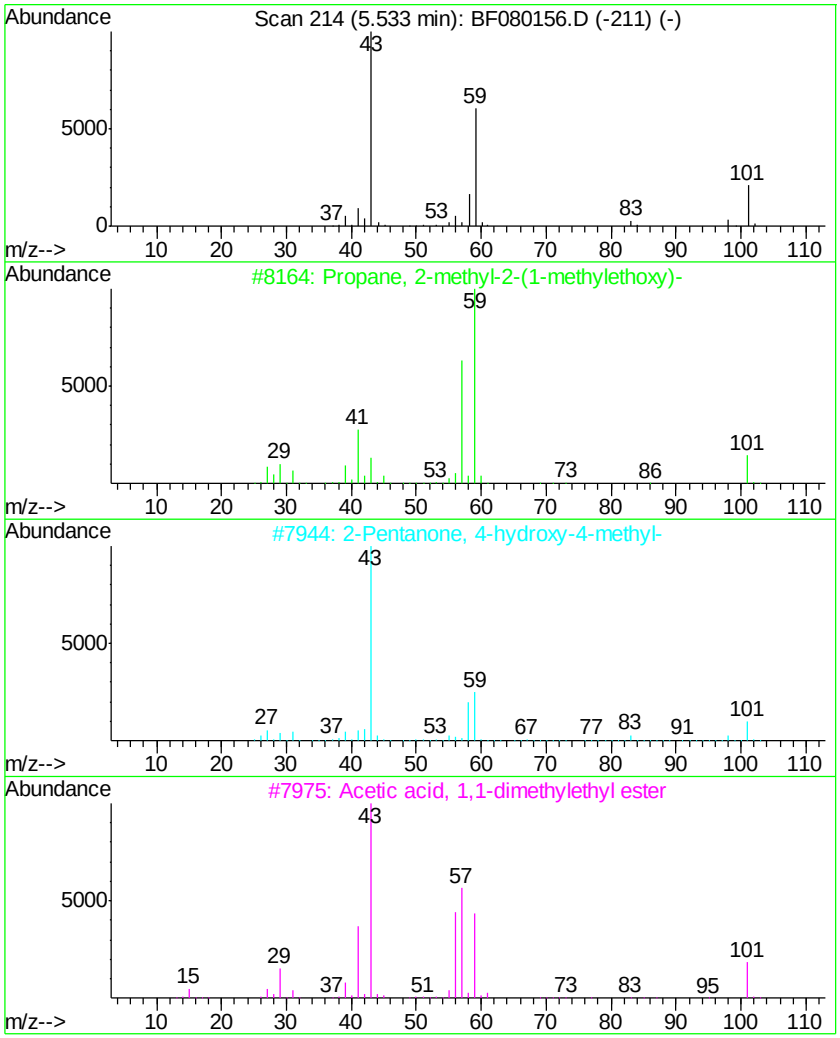
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

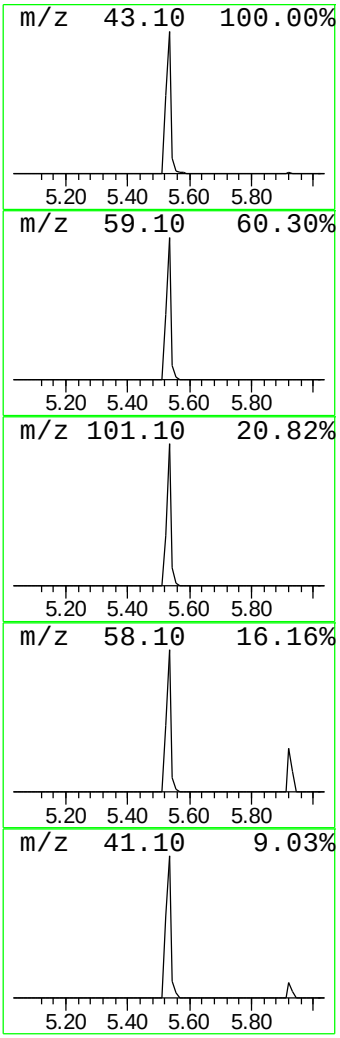
****

Peak Number 1 Propane, 2-methyl-2-(1-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	35.13 ng	345695	1,4-Dichlorobenzene-d4	7.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





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Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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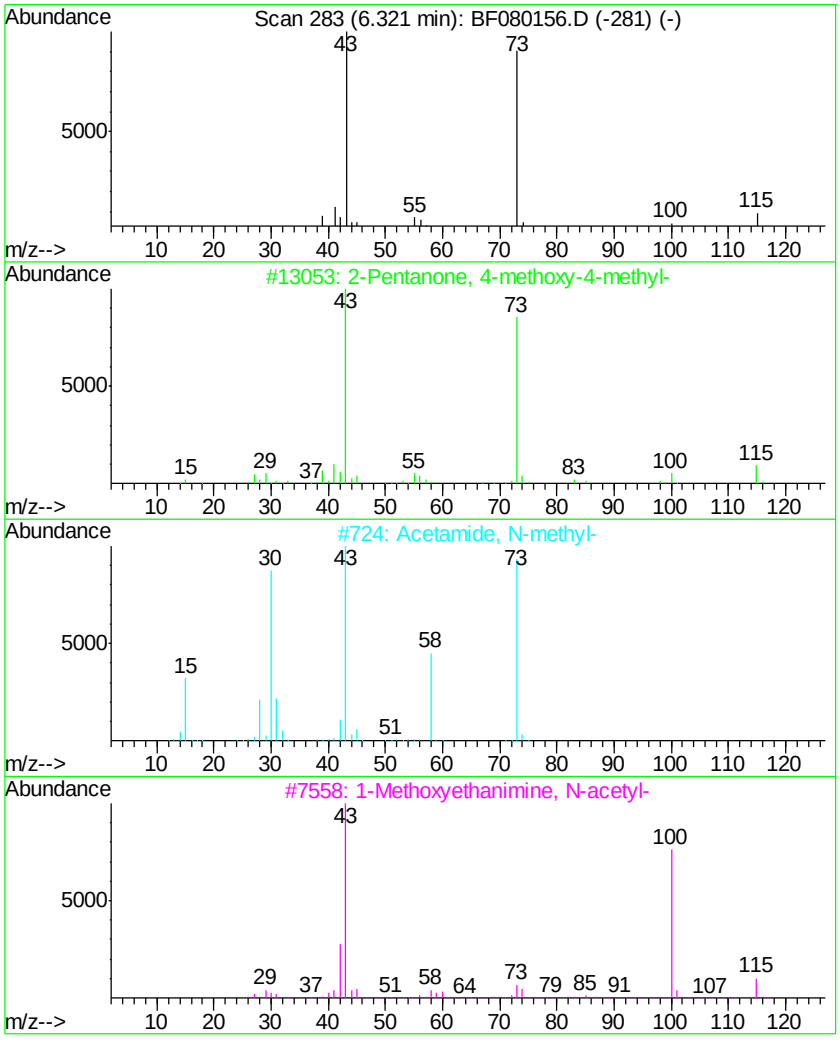
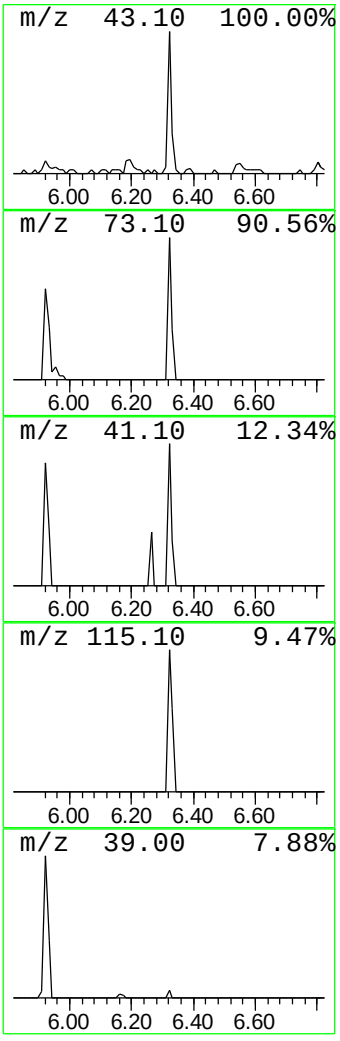
Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	2.81 ng	27653	1,4-Dichlorobenzene-d4	7.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	50
3	1-Methoxyethanimine, N-acetyl-	115	C5H9NO2	099028-43-0	9
4	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5	5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9





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Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
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TIC Library : C:\DATABASE\NIST02.L
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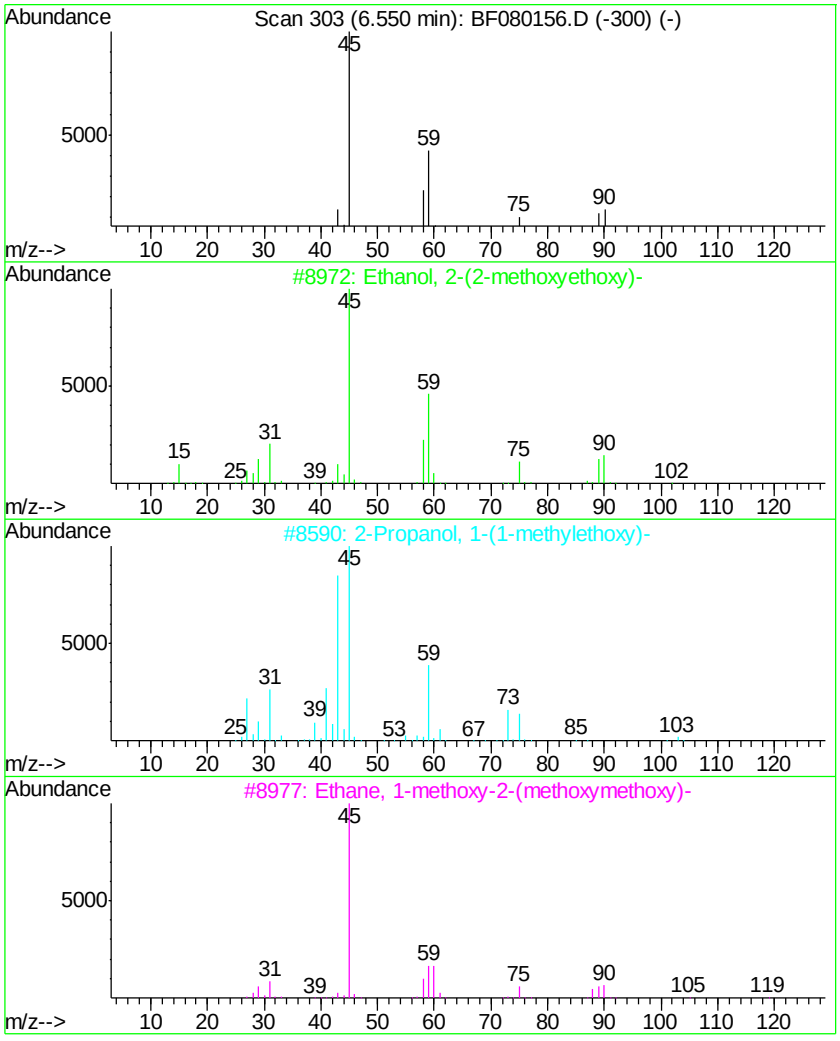
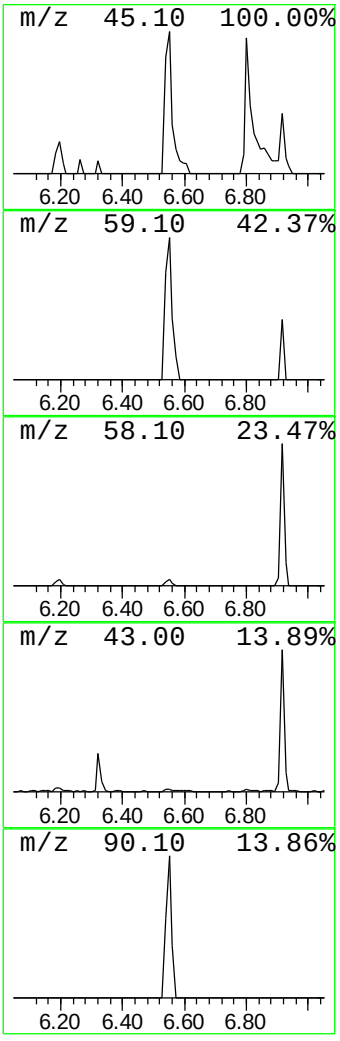
Peak Number 4 Ethanol, 2-(2-methoxyethoxy)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	2.19 ng	21567	1,4-Dichlorobenzene-d4	7.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	83
2	2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	9
3	Ethane, 1-methoxy-2-(methoxymeth...	120	C5H12O3	074498-88-7	9
4	Acetic acid, hydroxy-, methyl ester	90	C3H6O3	000096-35-5	7
5	Carbonic acid, dimethyl ester	90	C3H6O3	000616-38-6	5





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
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Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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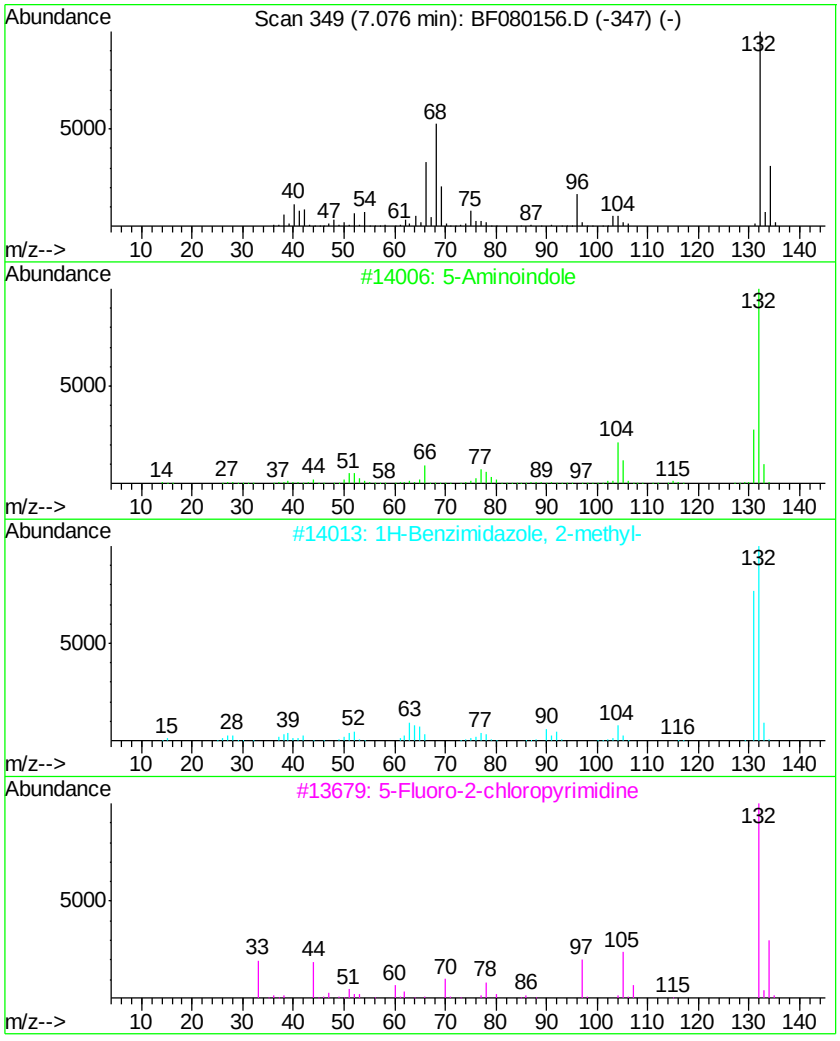
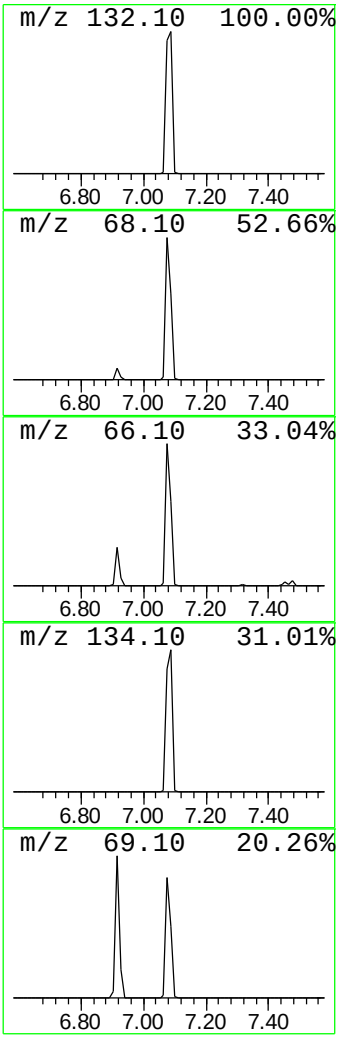
Peak Number 5 unknown7.08 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	107.74 ng	1060190	1,4-Dichlorobenzene-d4	7.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Aminoindole	132	C8H8N2	005192-03-0	22
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	11
5	(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10





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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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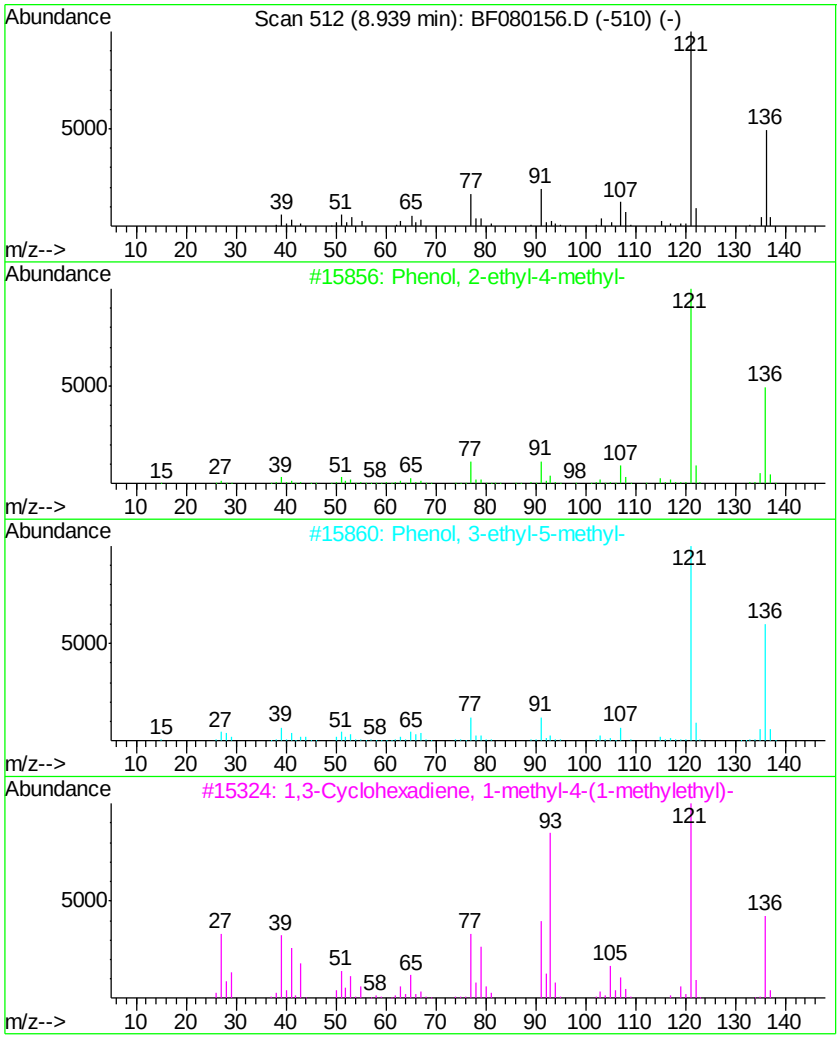
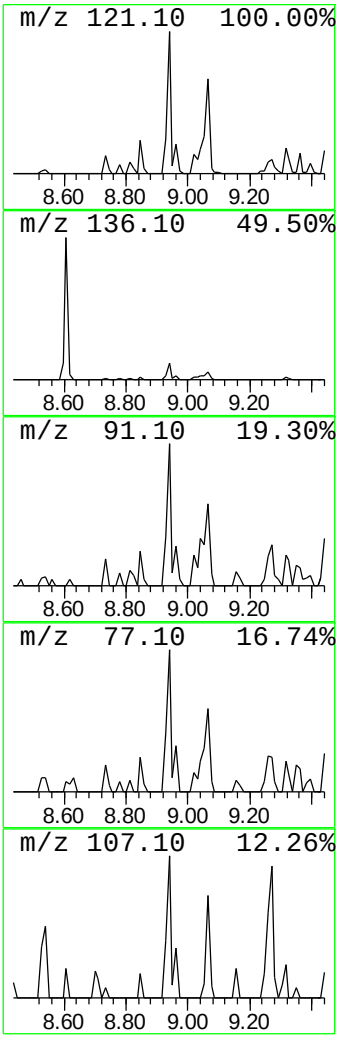
Peak Number 7 Phenol, 2-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.94	8.20 ng	109720	Naphthalene-d8	8.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-ethyl-4-methyl-	136	C9H12O	003855-26-3	91
2	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	91
3	1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	86
4	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	83
5	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	80





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Sample : G2699-01
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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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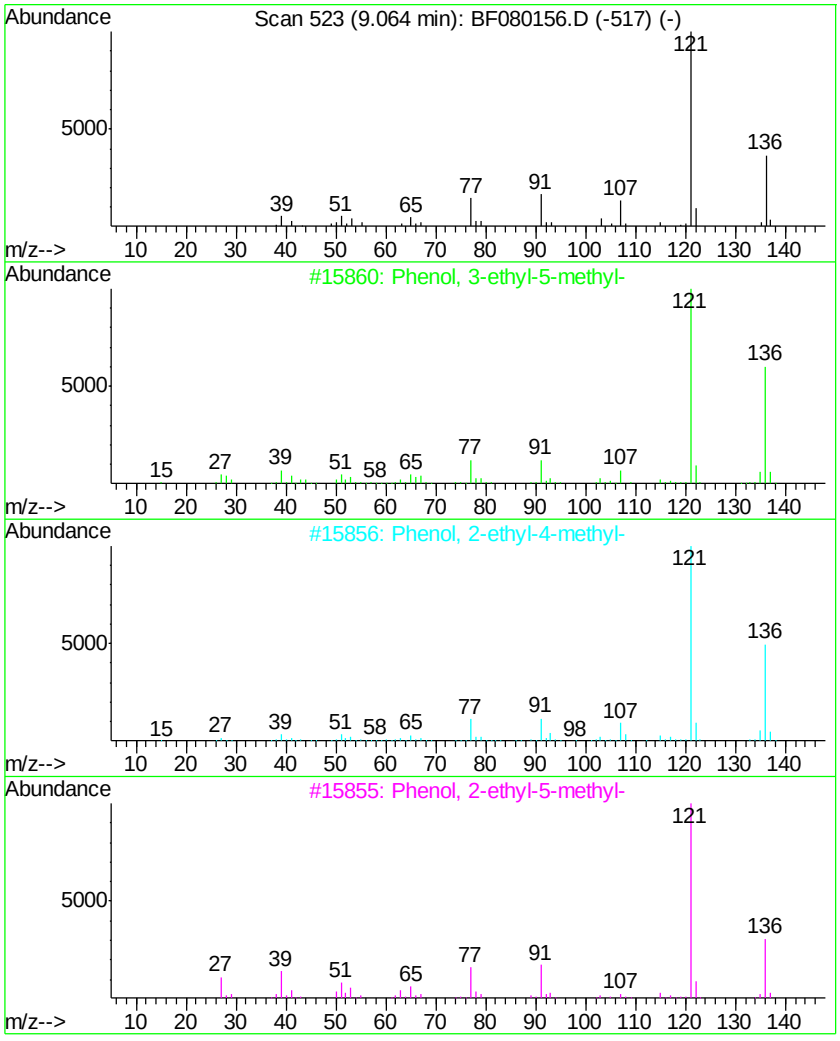
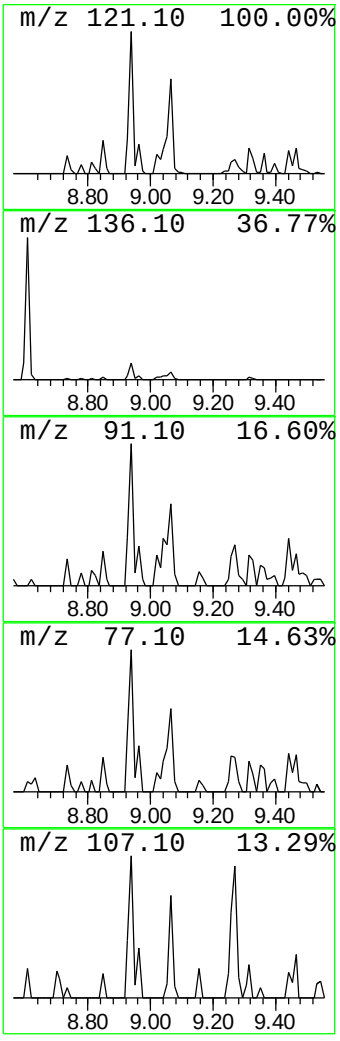
Peak Number 8 Phenol, 3-ethyl-5-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.06	8.02 ng	107364	Naphthalene-d8	8.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	91
2	Phenol, 2-ethyl-4-methyl-	136	C9H12O	003855-26-3	91
3	Phenol, 2-ethyl-5-methyl-	136	C9H12O	001687-61-2	90
4	2-Methylbicyclo[4.3.0]non-1(6)-ene	136	C10H16	060223-07-6	90
5	Phenol, 4-ethyl-3-methyl-	136	C9H12O	001123-94-0	87





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Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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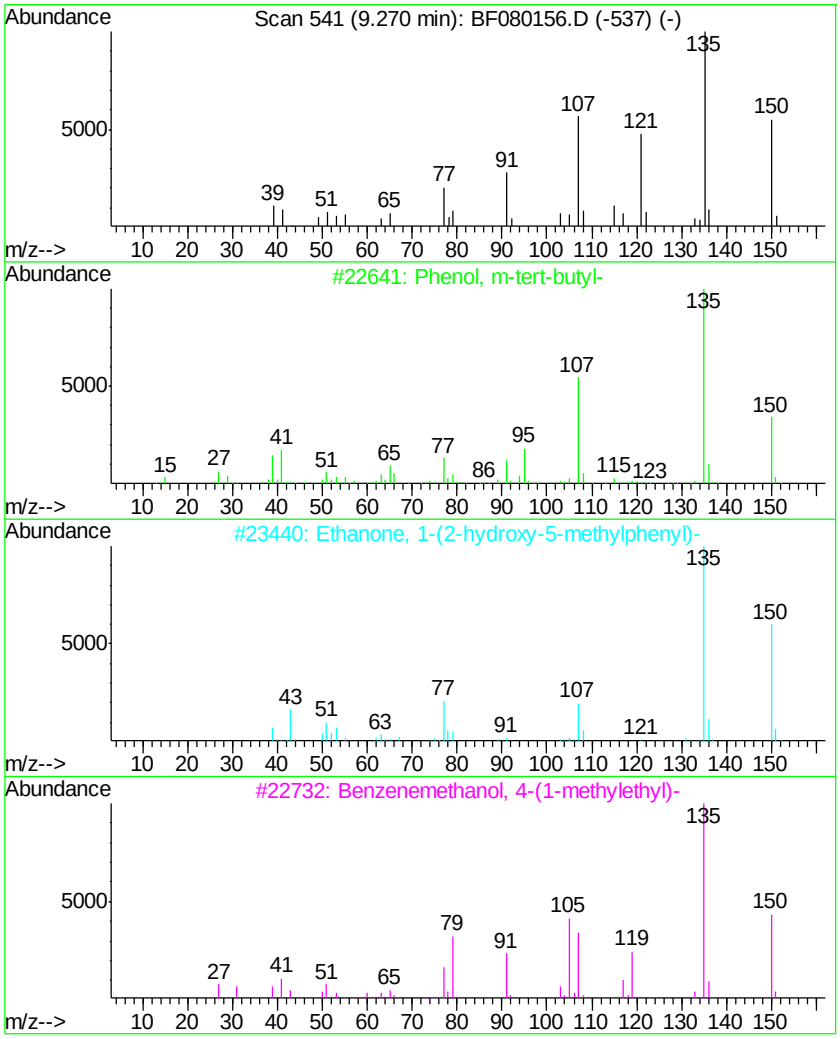
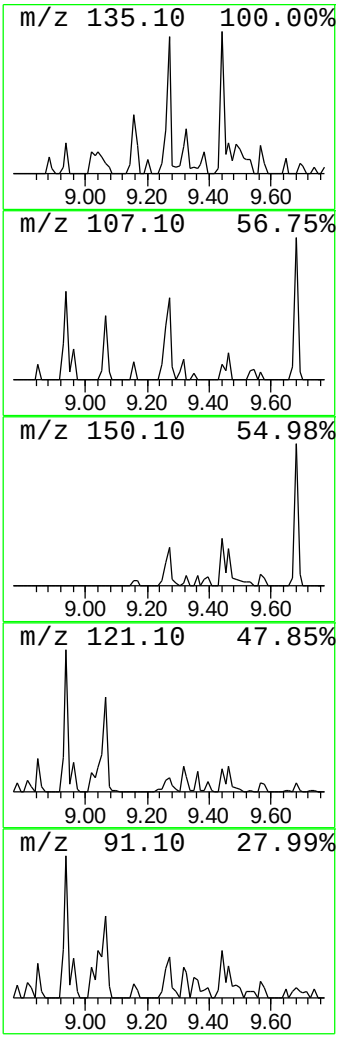
Peak Number 9 Phenol, m-tert-butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.27	3.65 ng	48908	Naphthalene-d8	8.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
2	Ethanone, 1-(2-hydroxy-5-methylp...	150	C9H10O2	001450-72-2	76
3	Benzenemethanol, 4-(1-methylethyl)-	150	C10H14O	000536-60-7	72
4	Ethanone, 1-(3-methoxyphenyl)-	150	C9H10O2	000586-37-8	72
5	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	72





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Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

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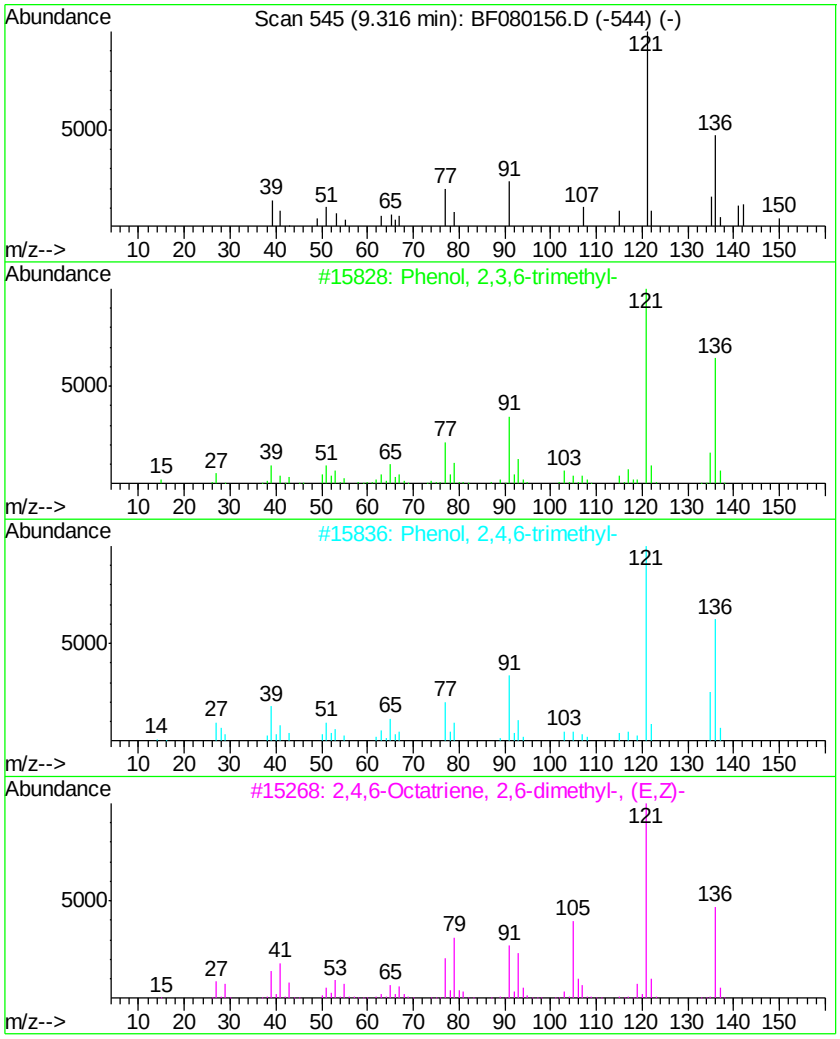
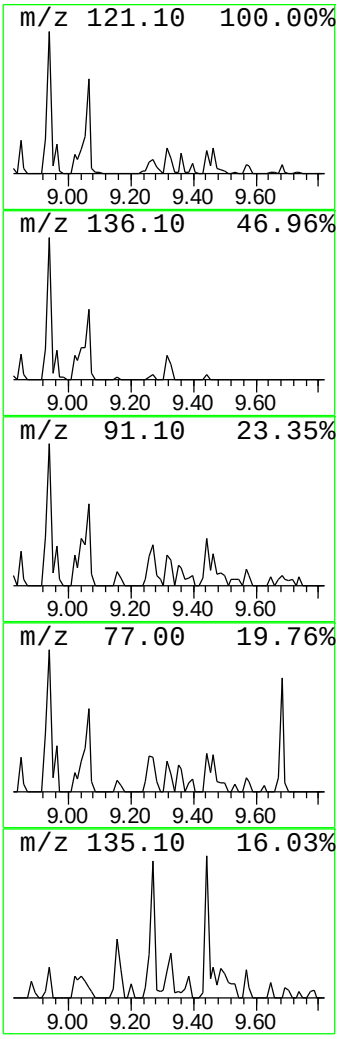
Peak Number 10 Phenol, 2,3,6-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.32	2.20 ng	29474	Naphthalene-d8	8.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	74
2	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	72
3	2,4,6-Octatriene, 2,6-dimethyl-,...	136	C10H16	007216-56-0	72
4	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	64
5	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	64





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080156.D
Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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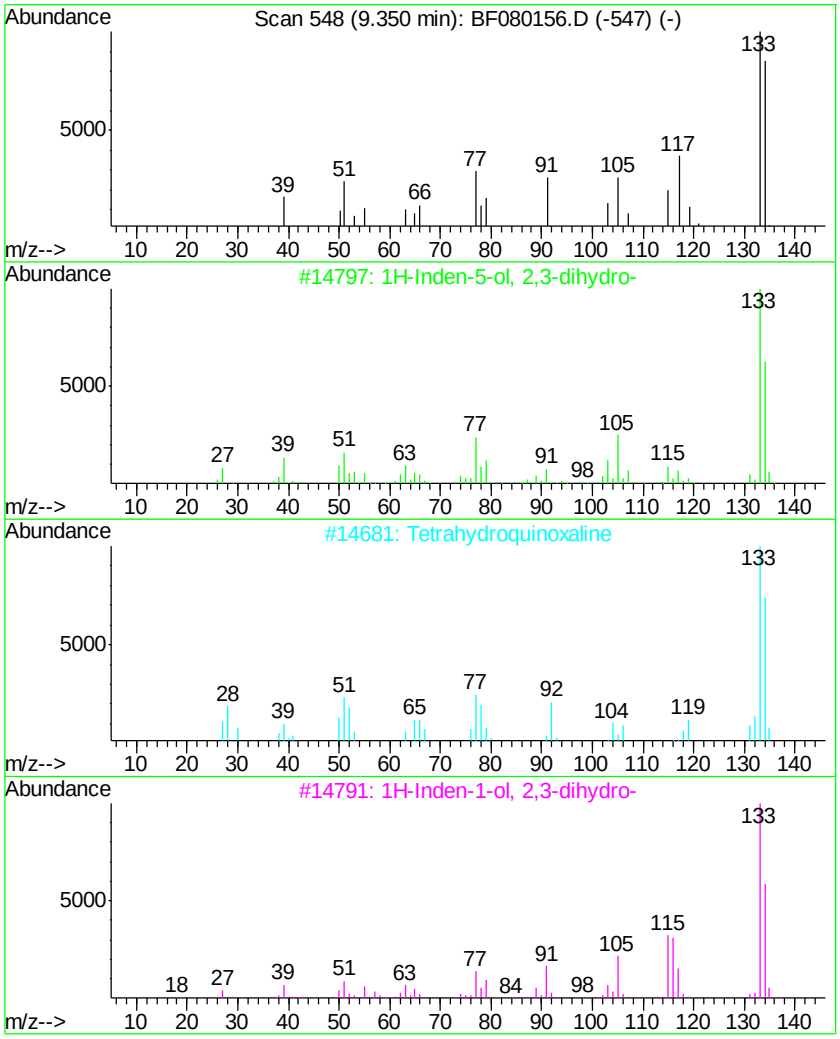
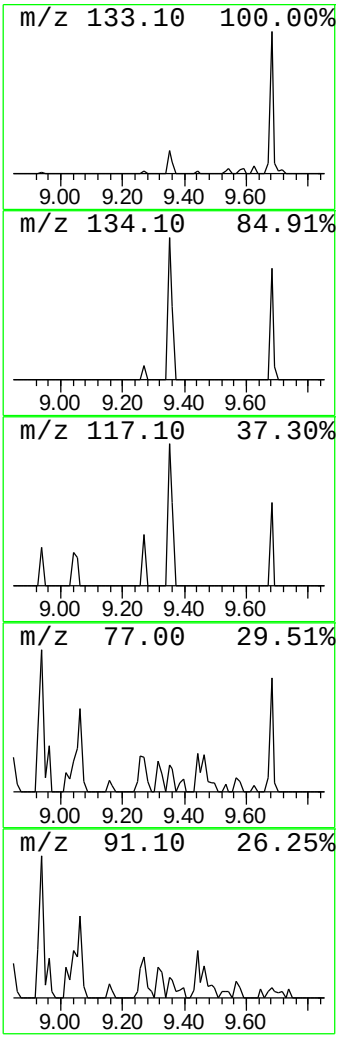
Peak Number 11 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.35	2.07 ng	27660	Naphthalene-d8	8.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	68
2	Tetrahydroquinoxaline	134	C8H10N2	003476-89-9	64
3	1H-Inden-1-ol, 2,3-dihydro-	134	C9H10O	006351-10-6	59
4	Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	59
5	Benzenemethanol, .alpha.-ethenyl-	134	C9H10O	004393-06-0	53





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Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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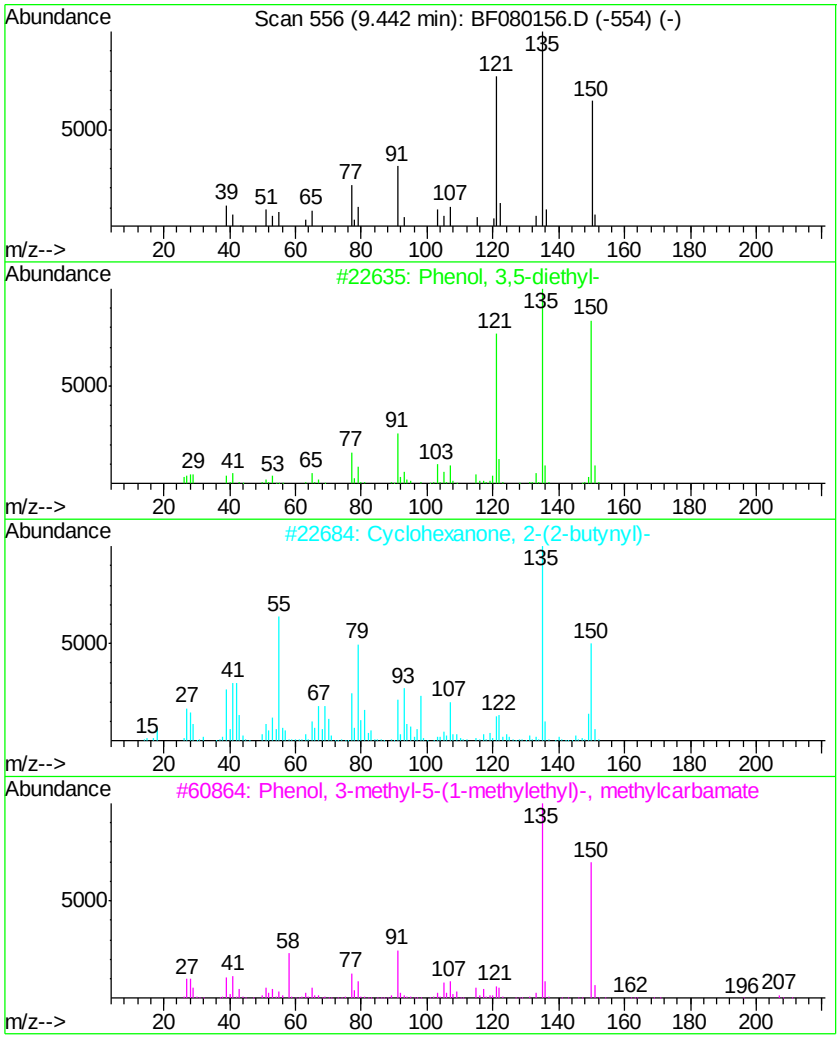
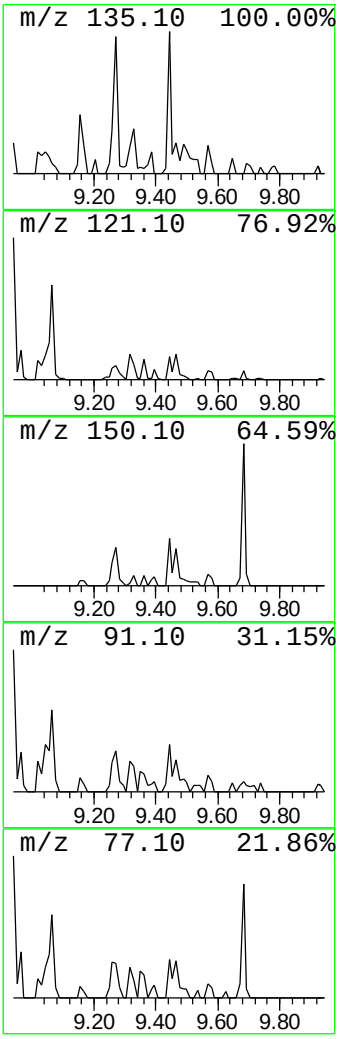
Peak Number 12 Phenol, 3,5-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	3.83 ng	51294	Naphthalene-d8	8.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-diethyl-	150	C10H14O	001197-34-8	91
2	Cyclohexanone, 2-(2-butynyl)-	150	C10H14O	054166-48-2	52
3	Phenol, 3-methyl-5-(1-methylethy...	207	C12H17NO2	002631-37-0	50
4	3-Methyl-4-isopropylphenol	150	C10H14O	003228-02-2	50
5	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	49





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Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
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TIC Library : C:\DATABASE\NIST02.L
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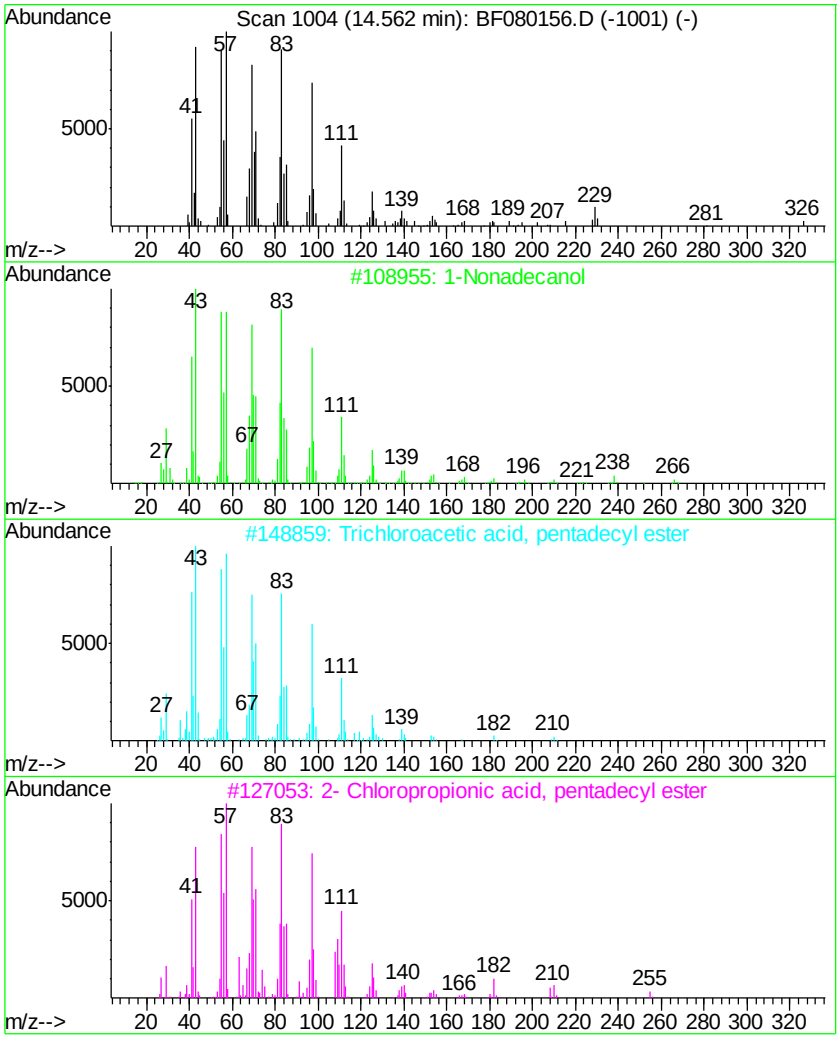
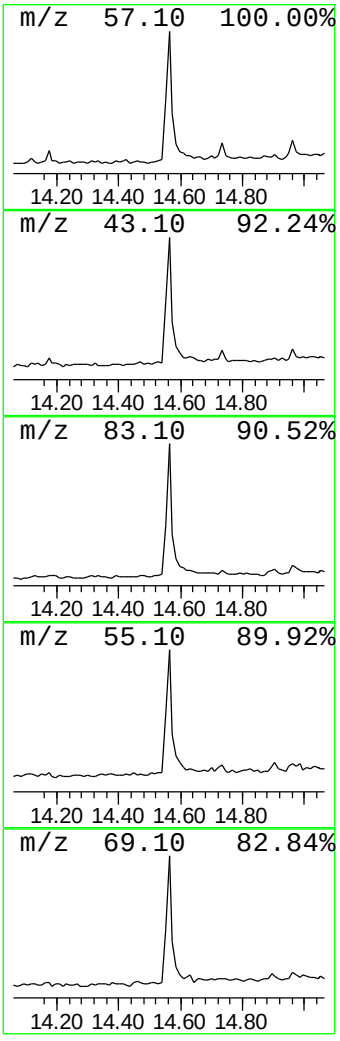
Peak Number 13 1-Nonadecanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	8.24 ng	246658	Chrysene-d12	14.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecanol	284	C19H40O	001454-84-8	95
2	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
3	2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	94
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5	1-Octadecanol	270	C18H38O	000112-92-5	93





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Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
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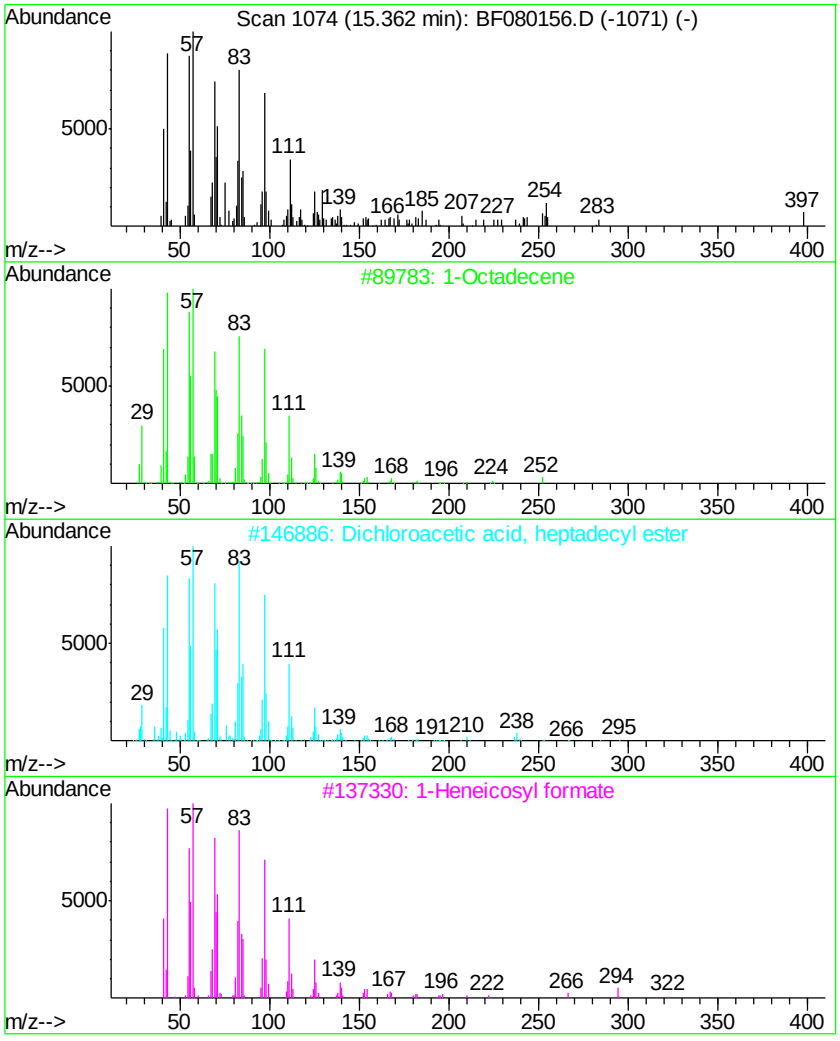
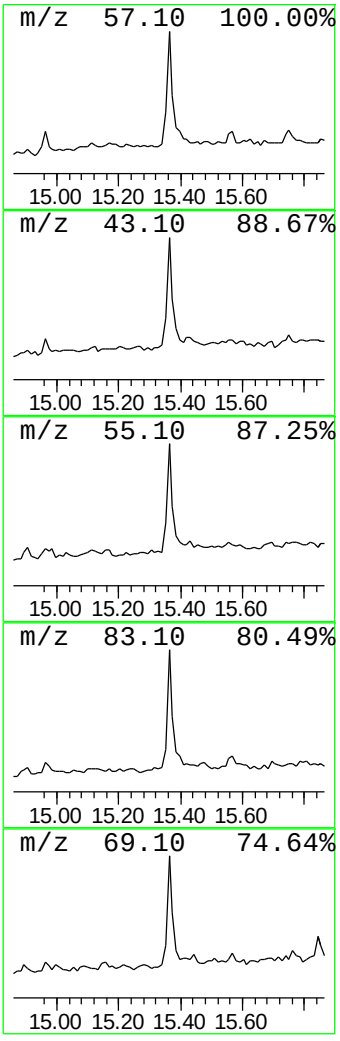
Peak Number 14 1-Octadecene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	6.43 ng	192438	Chrysene-d12	14.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	97
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94





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Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : LAR9819

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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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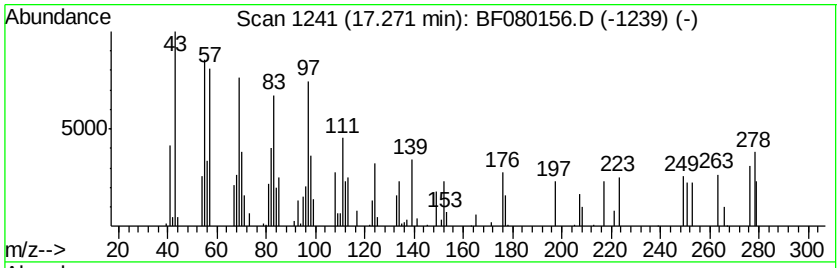
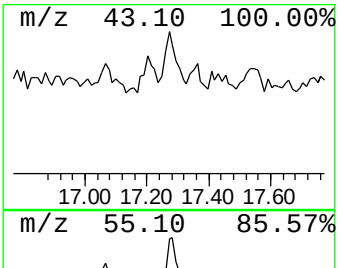
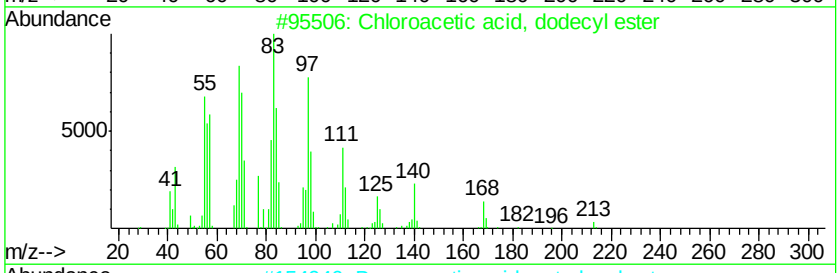
Peak Number 16 unknown17.27 Concentration Rank 12

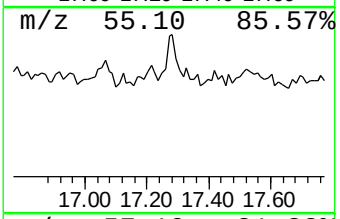
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.27	2.79 ng	61225	Perylene-d12	16.62

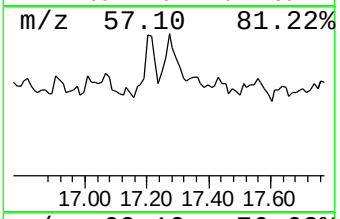
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chloroacetic acid, dodecyl ester	262	C14H27ClO2	006316-04-7	43
2	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	43
3	2-Cyclohexen-1-ol, 1-methyl-	112	C7H12O	023758-27-2	38
4	4-Methyl-E-9-octadecene	266	C19H38	1000130-87-1	30
5	Z-5-Nonadecene	266	C19H38	1000131-11-8	30

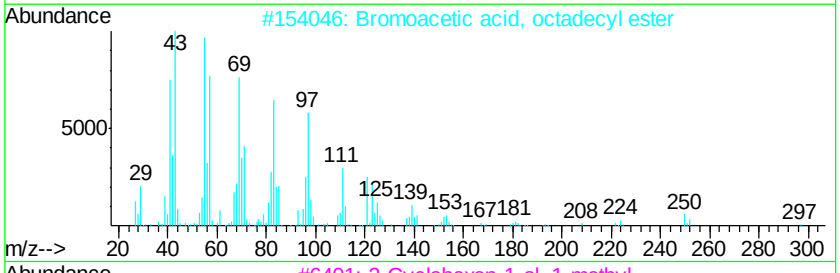


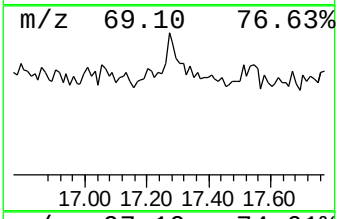


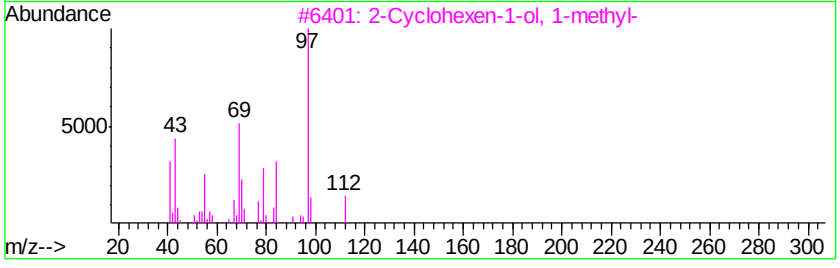


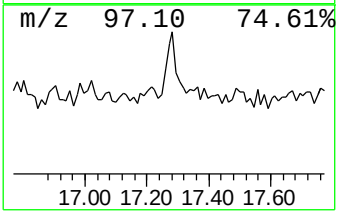












Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080156.D
Acq On : 25 Jun 2015 5:02
Operator : TP/IZ
Sample : G2699-01
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Instrument : BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Propane, 2-methyl...	5.53	35.1	ng	345695	1	7.32	196801	20.0
2-Pentanone, 4-me...	6.32	2.8	ng	27653	1	7.32	196801	20.0
Ethanol, 2-(2-met...	6.55	2.2	ng	21567	1	7.32	196801	20.0
unknown7.08	7.08	107.7	ng	1060190	1	7.32	196801	20.0
Phenol, 2-ethyl-4...	8.94	8.2	ng	109720	2	8.61	267720	20.0
Phenol, 3-ethyl-5...	9.06	8.0	ng	107364	2	8.61	267720	20.0
Phenol, m-tert-bu...	9.27	3.6	ng	48908	2	8.61	267720	20.0
Phenol, 2,3,6-tri...	9.32	2.2	ng	29474	2	8.61	267720	20.0
1H-Inden-5-ol, 2,...	9.35	2.1	ng	27660	2	8.61	267720	20.0
Phenol, 3,5-diethyl-	9.44	3.8	ng	51294	2	8.61	267720	20.0
1-Nonadecanol	14.56	8.2	ng	246658	5	14.78	598852	20.0
1-Octadecene	15.36	6.4	ng	192438	5	14.78	598852	20.0
unknown17.27	17.27	2.8	ng	61225	6	16.62	438176	20.0