LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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	3.144	3	5	8	rVB	1516302	2472535	100.00%	17.425%
2	3.738	54	57	60	rBV	86847	126434	5.11%	0.891%
3	4.824	150	152	155	rBV	67619	82932	3.35%	0.584%
4	5.430	204	205	208	rVV	46676	43330	1.75%	0.305%
5	5.499	208	211	215	rVB	41832	47278	1.91%	0.333%
6	5.887	243	245	248	rBV	385694	488849	19.77%	3.445%
7	6.139	265	267	270	rBV	45531	45557	1.84%	0.321%
8	6.367	286	287	289	rVB	20974	24816	1.00%	0.175%
9	6.882	330	332	335	rBV	249643	326317	13.20%	2.300%
10	6.973	338	340	342	rVB	55762	43620	1.76%	0.307%
11	7.053	344	347	349	rBV	925253	888167	35.92%	6.259%
12	7.282	365	367	369	rBV	242395	210629	8.52%	1.484%
13	7.339	369	372	374	rVB	55758	50128	2.03%	0.353%
14	7.442	379	381	385	rVV	994246	904011	36.56%	6.371%
15	7.522	385	388	394	rVB2	64404	123377	4.99%	0.870%
16	7.796	407	412	414	rBV	537514	1076029	43.52%	7.583%
17	7.842	414	416	426	rVV	848302	910249	36.81%	6.415%
18	8.127	438	441	445	rVB2	44706	76678	3.10%	0.540%
19	8.310	456	457	459	rVB2	46537	34030	1.38%	0.240%
20	8.436	466	468	473	rVB5	9462	26825	1.08%	0.189%
21	8.573	478	480	483	rBV	358527	290056	11.73%	2.044%
22	9.110	526	527	531	rVB	33225	36668	1.48%	0.258%
23	9.522	562	563	567	rVB	45495	45263	1.83%	0.319%
24	9.648	572	574	576	rVB	1368739	1245253	50.36%	8.776%
25	9.819	587	589	596	rBV	125195	128437	5.19%	0.905%
26	10.345	633	635	638	rVB	411909	342107	13.84%	2.411%
27	10.951	686	688	690	rBV	634043	494284	19.99%	3.483%
28	11.133	702	704	707	rBV2	600592	767250	31.03%	5.407%
29	11.853	764	767	769	rBV	294647	337065	13.63%	2.375%
30	12.048	782	784	787	rBV	262242	218996	8.86%	1.543%
31	13.522	910	913	915	rBV	1498992	1536610	62.15%	10.829%
32	14.768	1019	1022	1024	rVB	313488	378164	15.29%	2.665%
33	16.597	1179	1182	1189	rVB	264350	367365	14.86%	2.589%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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

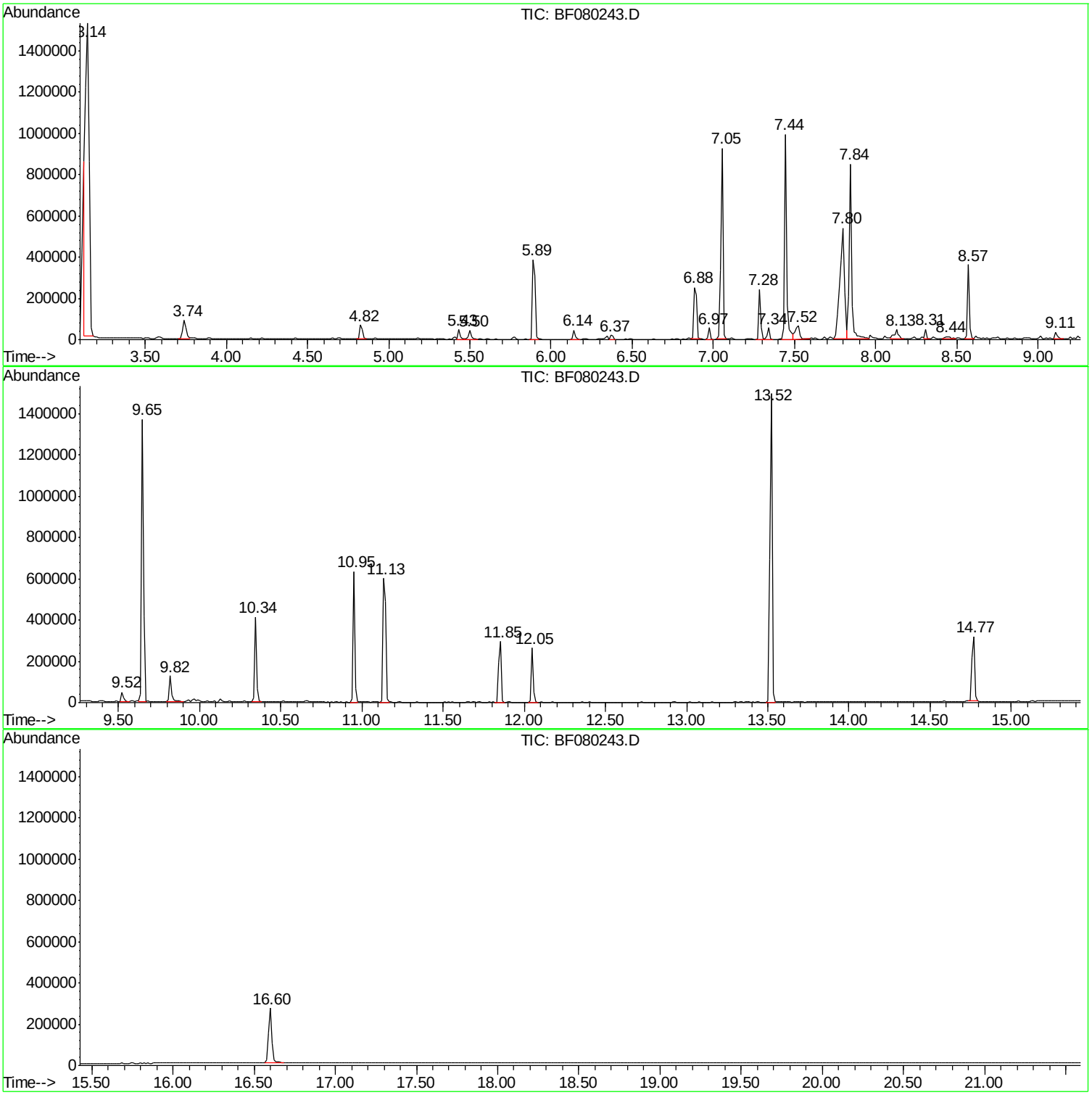
Sum of corrected areas: 14189309

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-1

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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

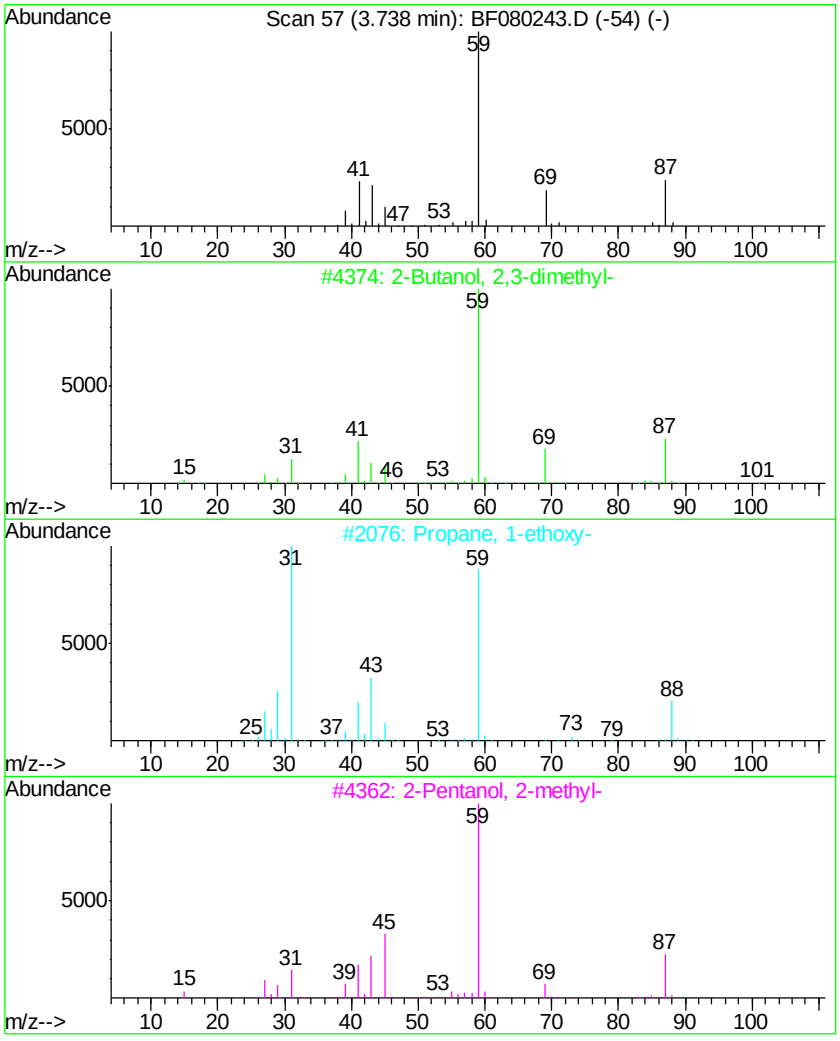
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

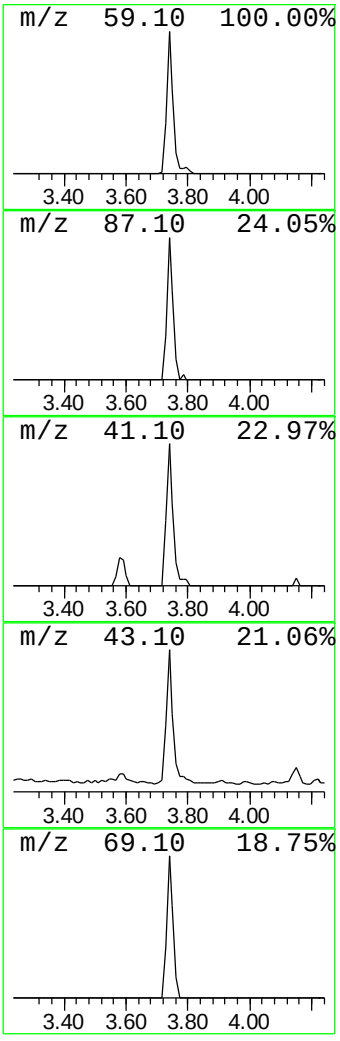
****

Peak Number 1 2-Butanol, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.74	12.01 ng	126434	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
2	Propane, 1-ethoxy-	88	C5H12O	000628-32-0	53
3	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	50
4	3-Heptanol	116	C7H16O	000589-82-2	40
5	3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	40





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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

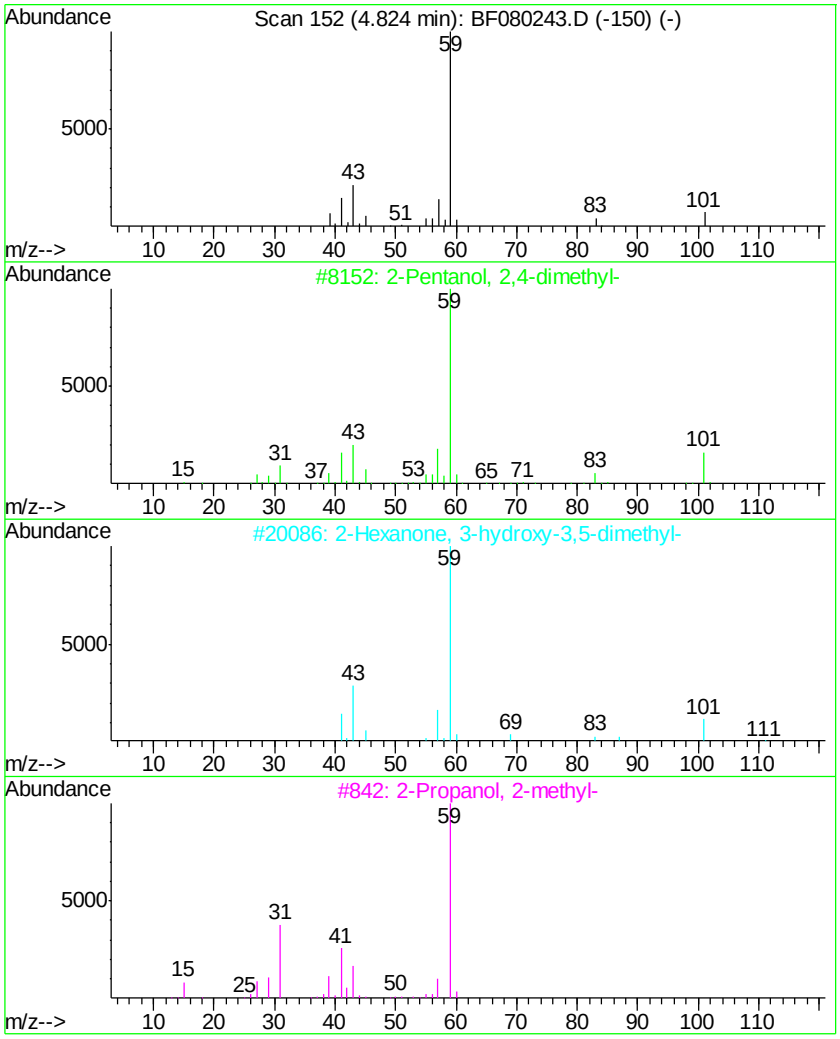
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

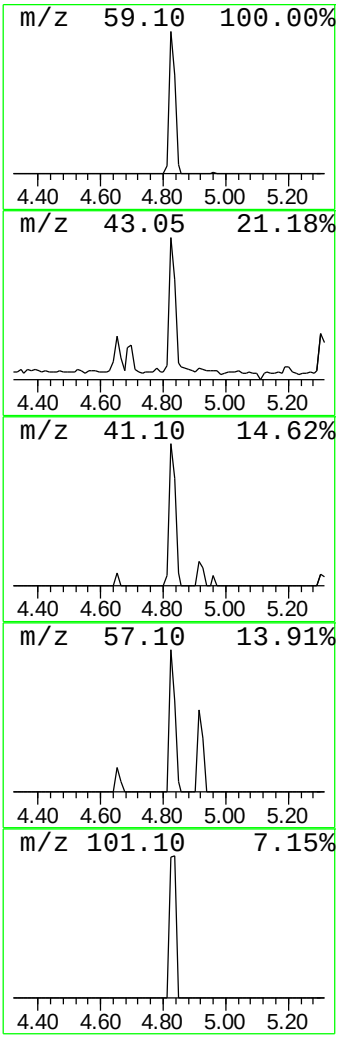
****

Peak Number 2 2-Pentanol, 2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.82	7.87 ng	82932	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	83
2	2-Hexanone, 3-hydroxy-3,5-dimethyl-	144	C8H16O2	006321-14-8	78
3	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	50
4	3-Pentanol	88	C5H12O	000584-02-1	50
5	2-Pentanol, 2,3-dimethyl-	116	C7H16O	004911-70-0	40





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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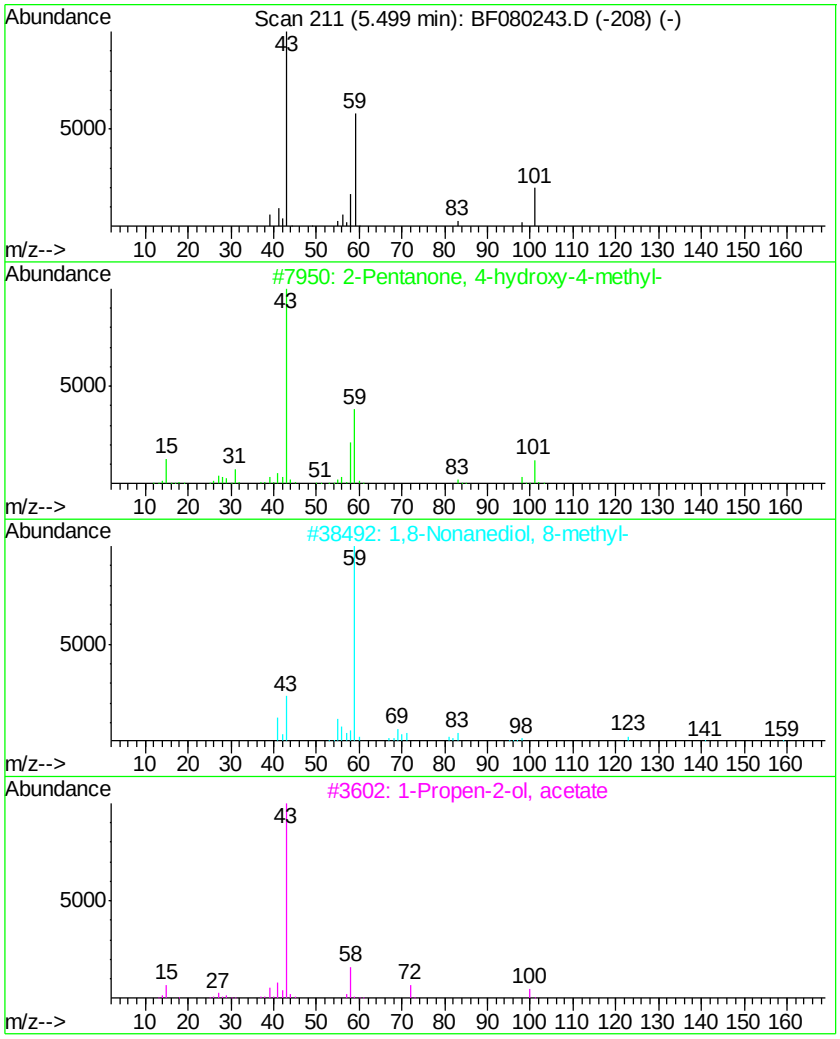
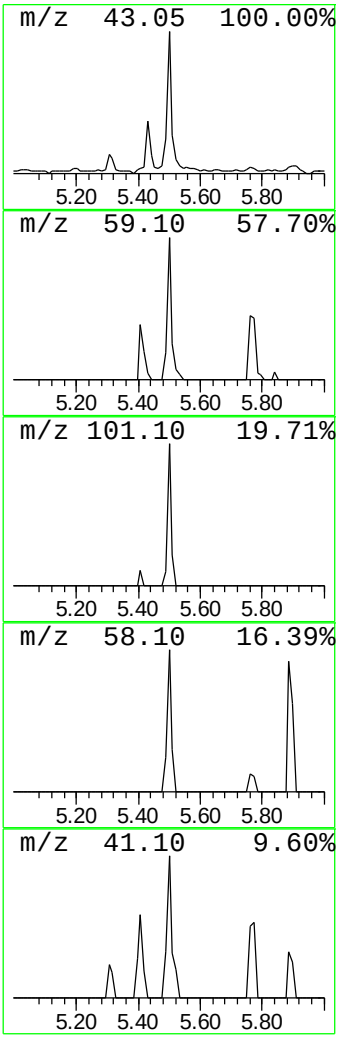
Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.50	4.49 ng	47278	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	5





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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

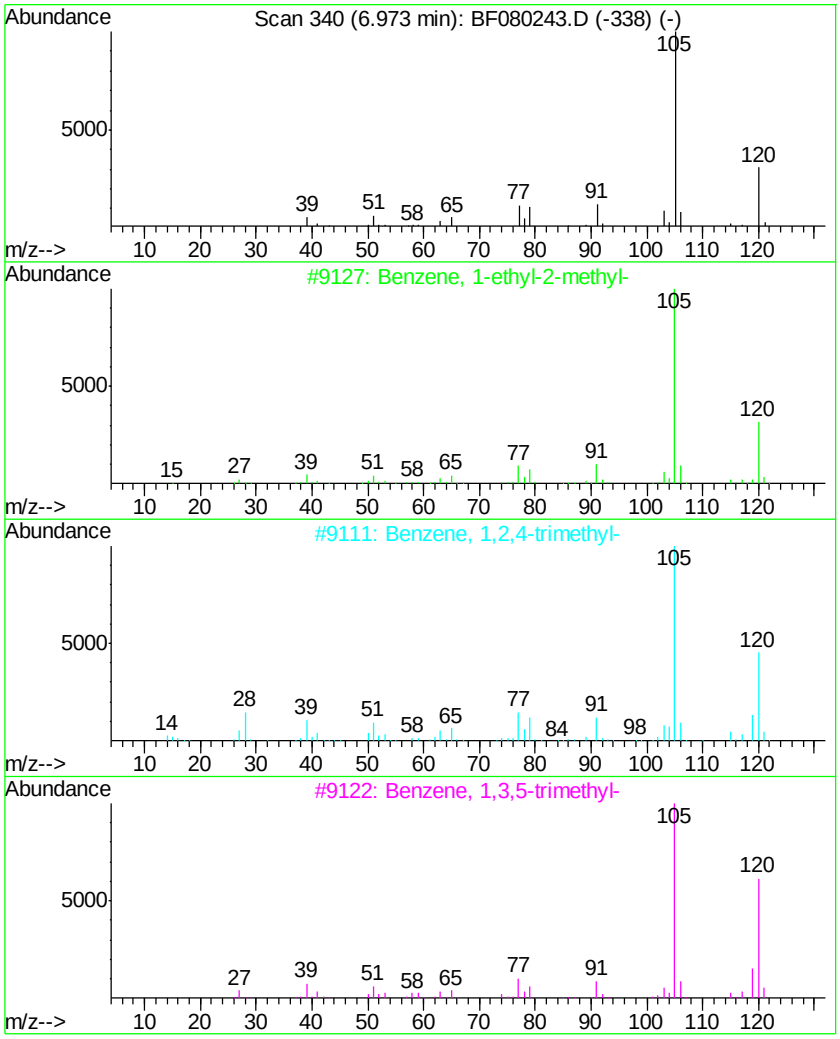
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

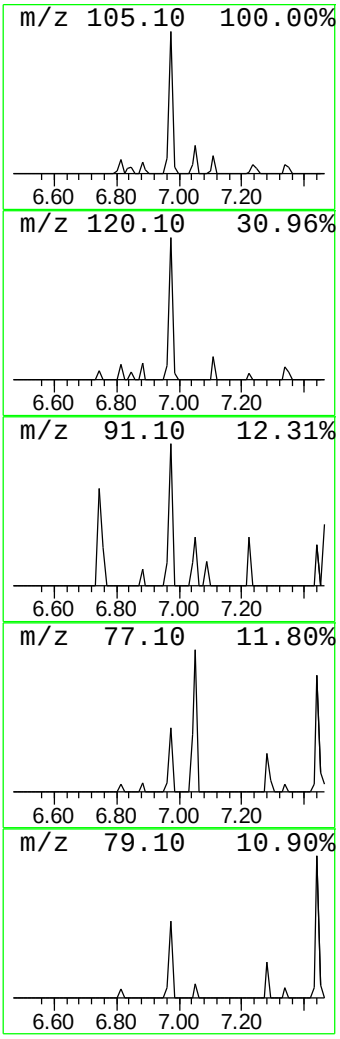
****

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.97	4.14 ng	43620	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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

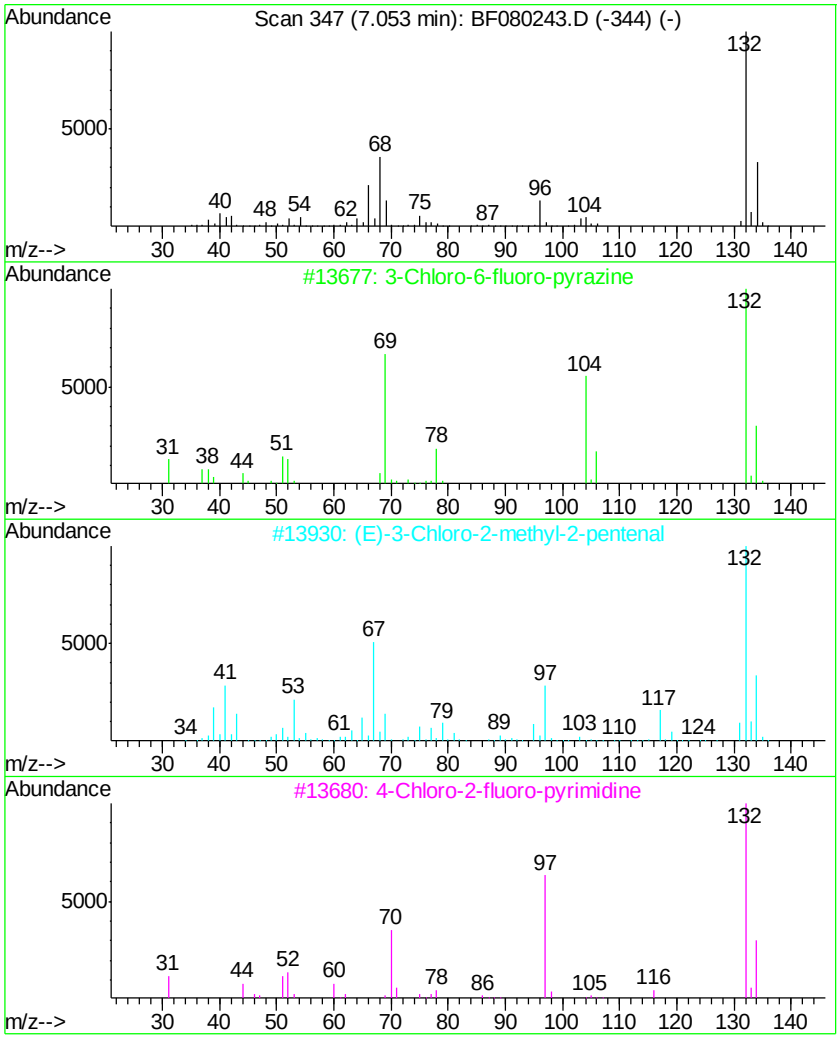
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

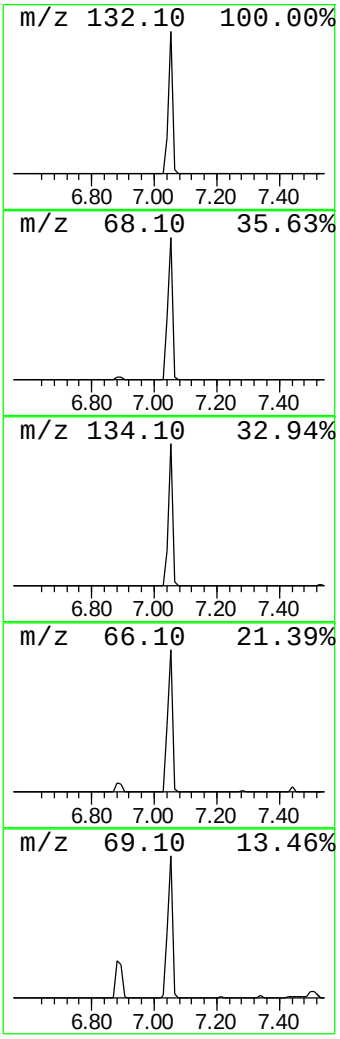
****

Peak Number 8 unknown7.05 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	84.33 ng	888167	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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

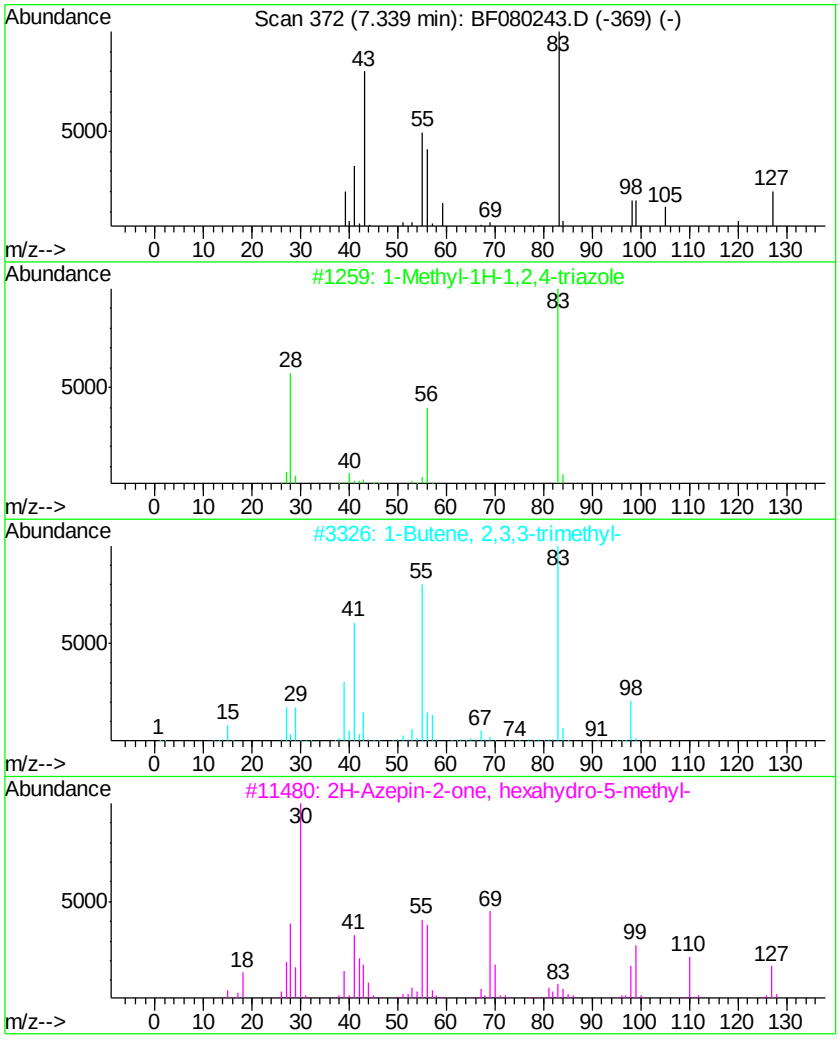
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

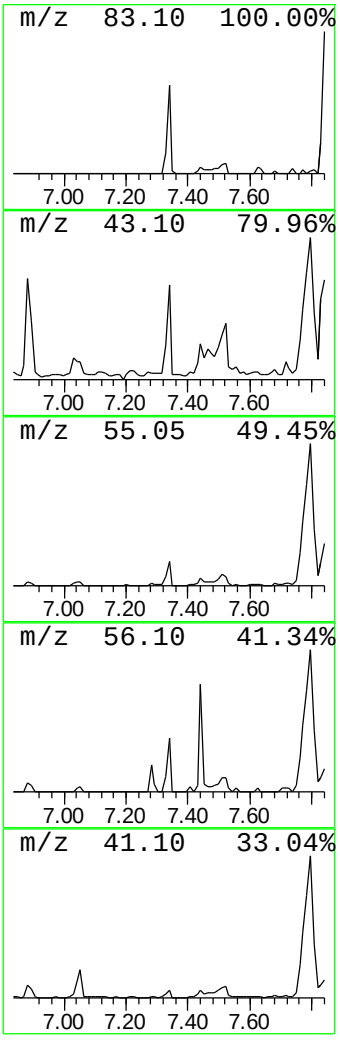
****

Peak Number 9 unknown7.34 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.34	4.76 ng	50128	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	40
2	1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	37
3	2H-Azepin-2-one, hexahydro-5-met...	127	C7H13NO	002210-07-3	27
4	3-Methyl-4-hexyn-3-ol	112	C7H12O	006320-68-9	25
5	1,5-Heptadien-4-one, 3,3,6-trime...	152	C10H16O	000546-49-6	25





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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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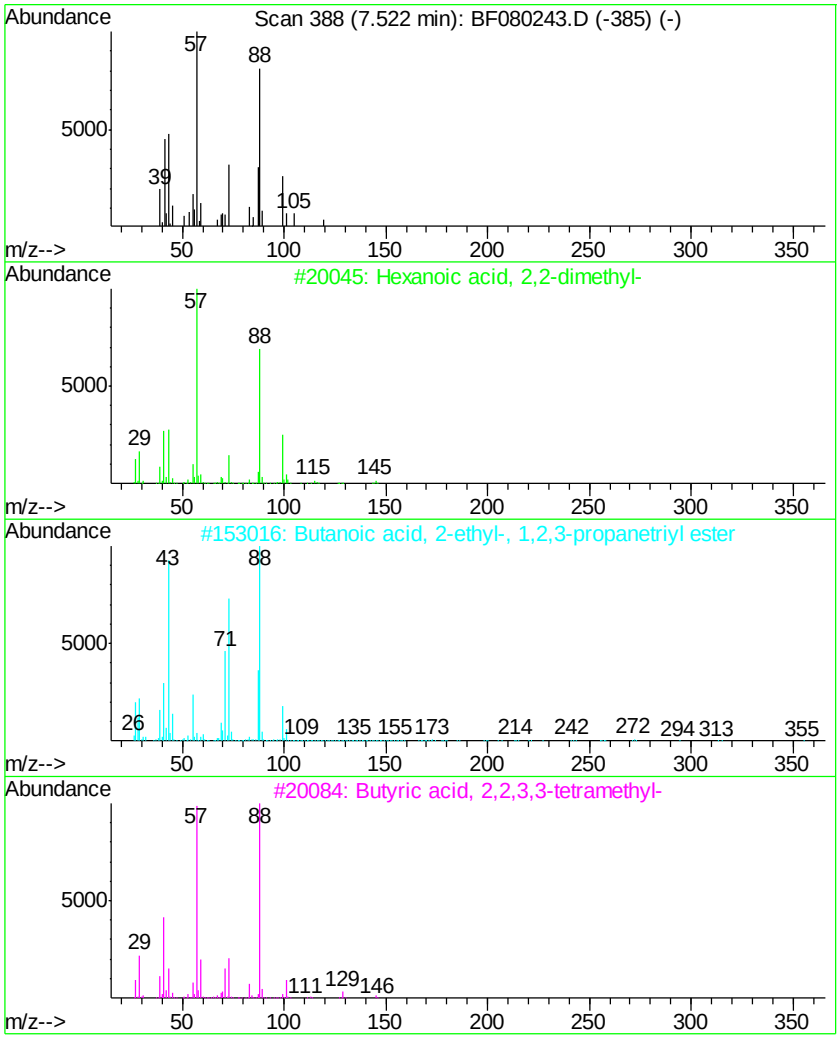
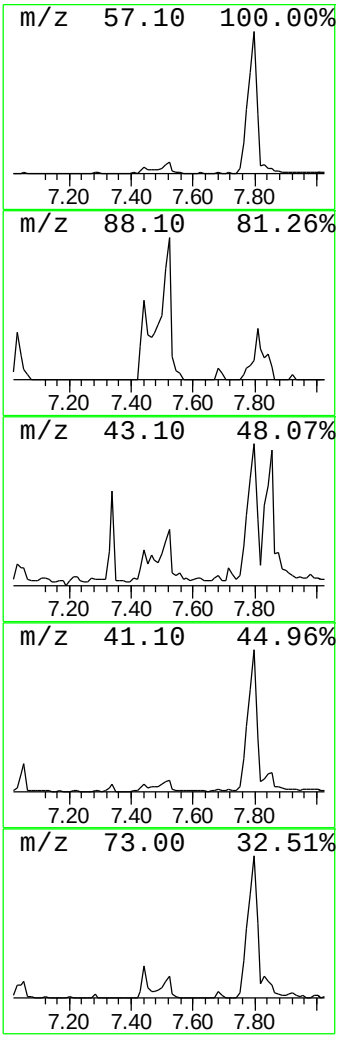
Peak Number 11 unknown7.52 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	11.72 ng	123377	1,4-Dichlorobenzene-d4	7.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2,2-dimethyl-	144	C8H16O2	000813-72-9	38
2	Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	36
3	Butyric acid, 2,2,3,3-tetramethyl-	144	C8H16O2	030407-41-1	33
4	Methyl 2-methylhexanoate	144	C8H16O2	002177-81-3	28
5	Cyclohexaneacetic acid, .alpha.-...	170	C10H18O2	006051-00-9	25





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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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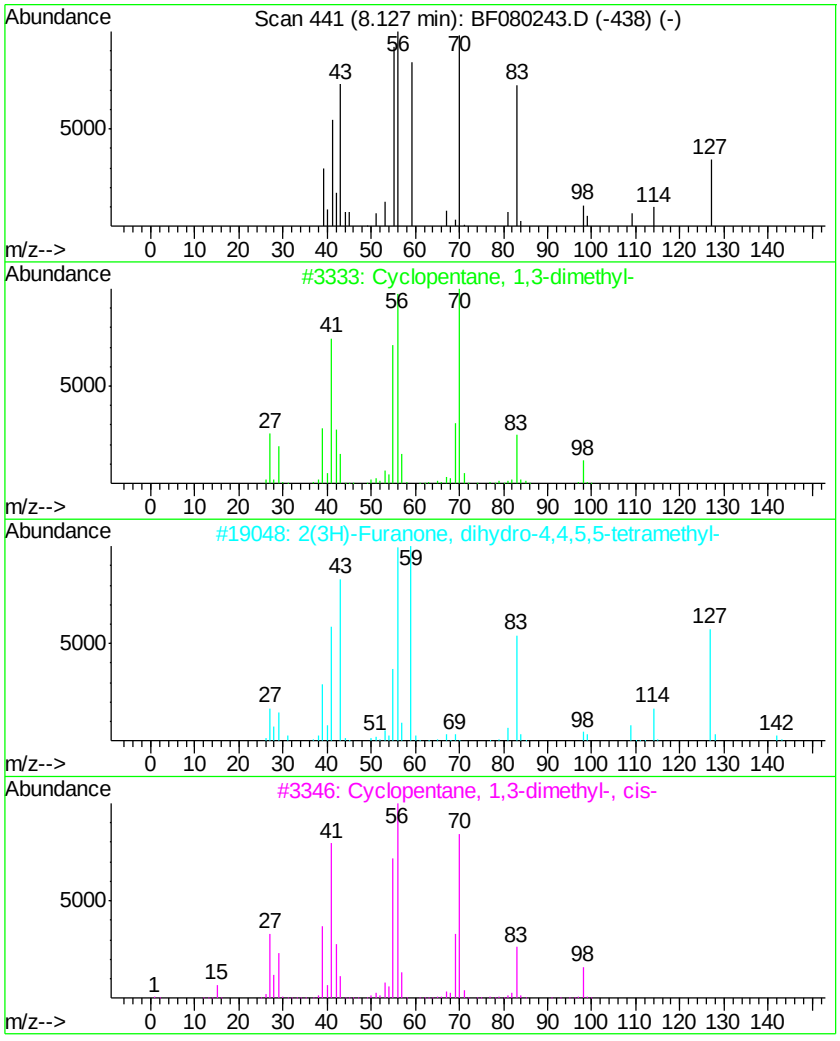
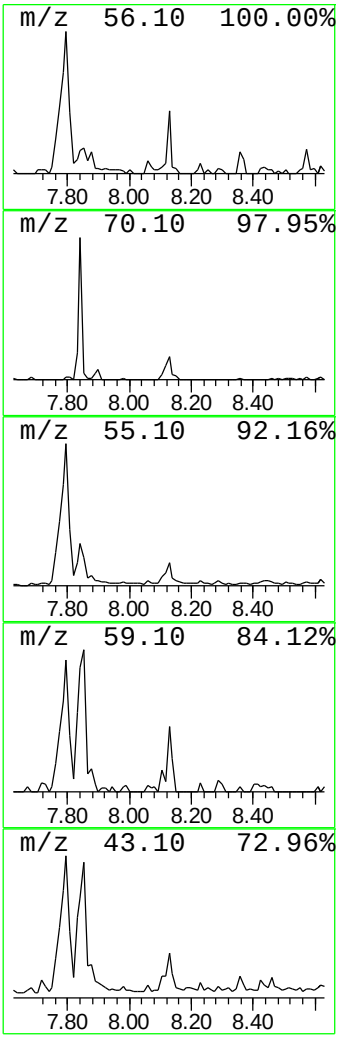
Peak Number 12 unknown8.13 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.13	5.29 ng	76678	Naphthalene-d8	8.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	43
2	2(3H)-Furanone, dihydro-4,4,5,5-...	142	C8H14O2	016466-24-3	43
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	43
4	Octane, 3-chloro-	148	C8H17Cl	001117-79-9	38
5	Cycloheptane	98	C7H14	000291-64-5	37





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF063015\
Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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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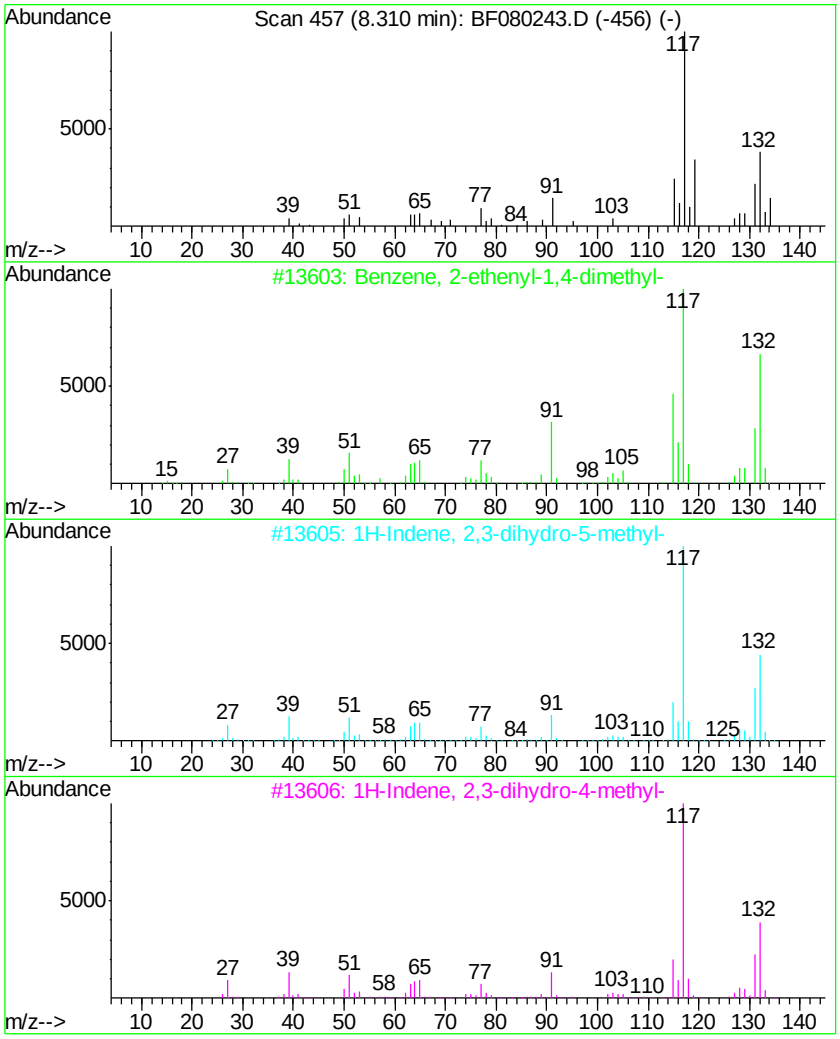
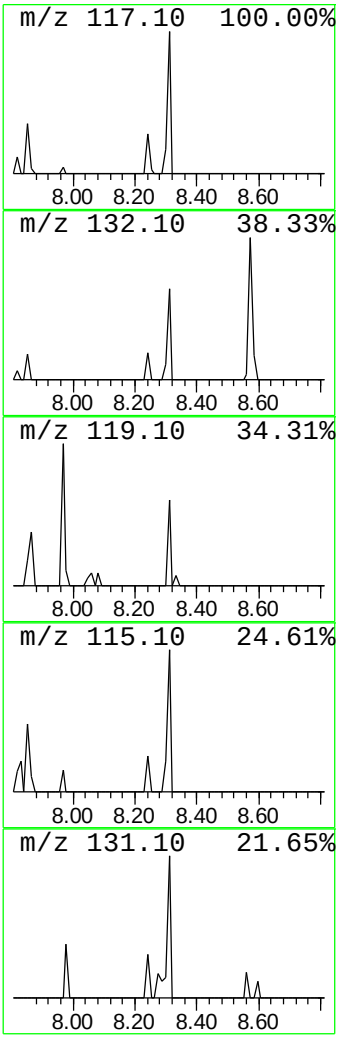
Peak Number 13 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	2.35 ng	34030	Naphthalene-d8	8.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76





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Data File : BF080243.D
Acq On : 30 Jun 2015 18:32
Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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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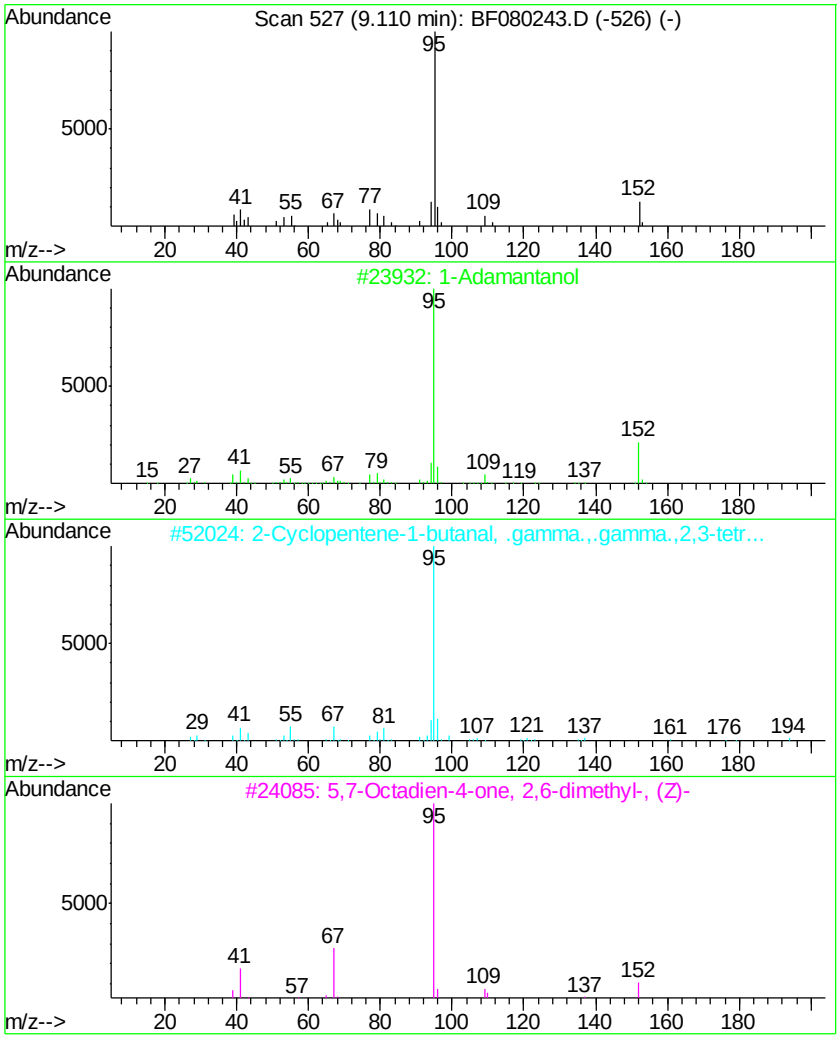
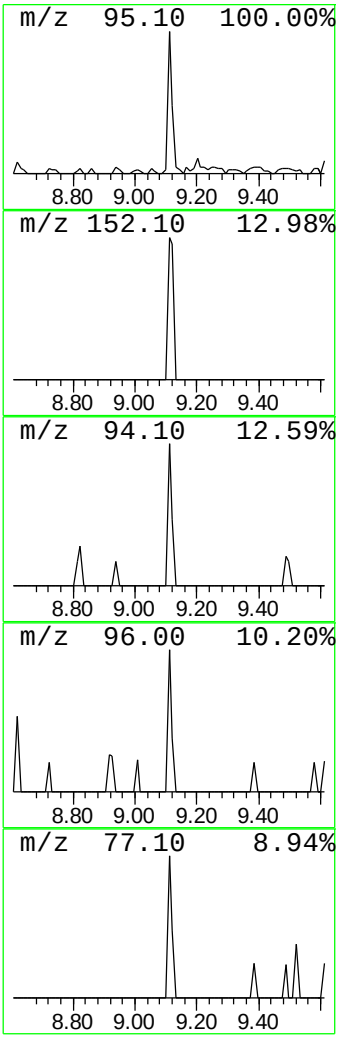
Peak Number 14 1-Adamantanol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.11	2.53 ng	36668	Naphthalene-d8	8.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Adamantanol	152	C10H16O	000768-95-6	91
2	2-Cyclopentene-1-butanal, .gamma...	194	C13H22O	060714-25-2	64
3	5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	003588-18-9	64
4	4-Pyridinol	95	C5H5NO	000626-64-2	58
5	Cyclohexanemethanol, 4-(1-methyl...	156	C10H20O	013828-37-0	39





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Sample : G2837-01
Misc :
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Instrument : BNA_F
ClientSampleId : MW-1

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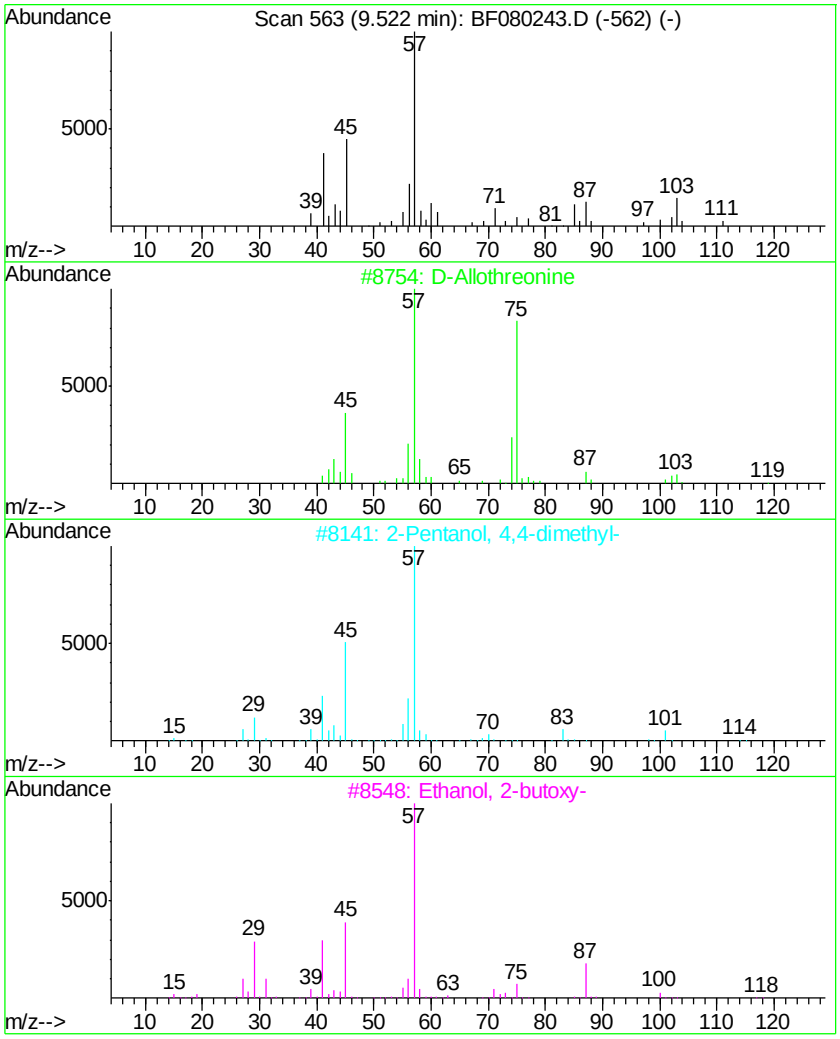
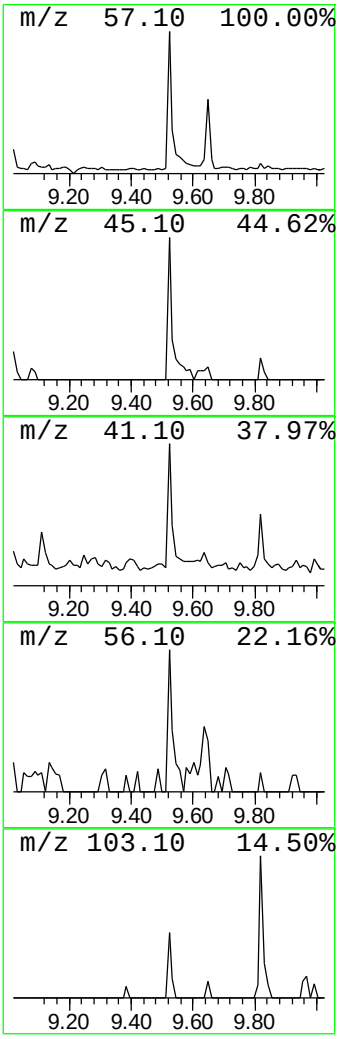
Peak Number 15 unknown9.52 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.52	2.65 ng	45263	Acenaphthene-d10	10.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	D-Allothreonine	119	C4H9NO3	024830-94-2	45
2	2-Pentanol, 4,4-dimethyl-	116	C7H16O	006144-93-0	43
3	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	40
4	Propene, 3-tert-butoxy-2-(methox...	158	C9H18O2	023230-86-6	38
5	dl-Threonine	119	C4H9NO3	000080-68-2	37





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Operator : TP/IZ
Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
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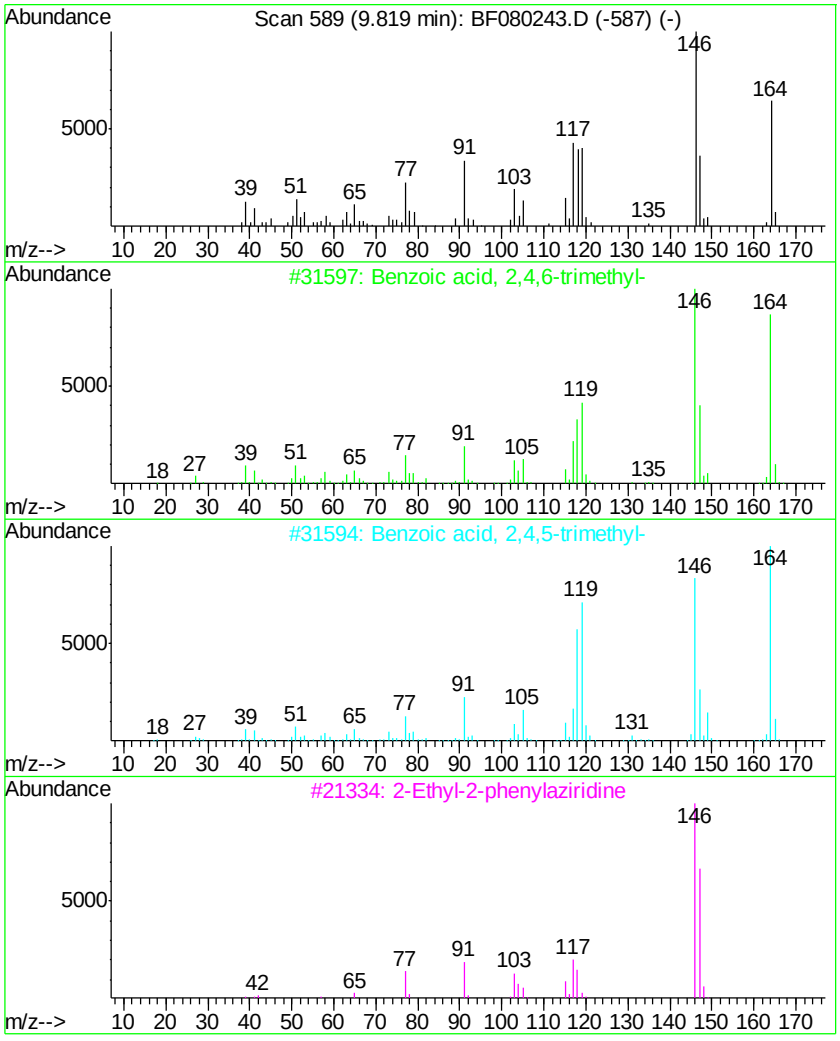
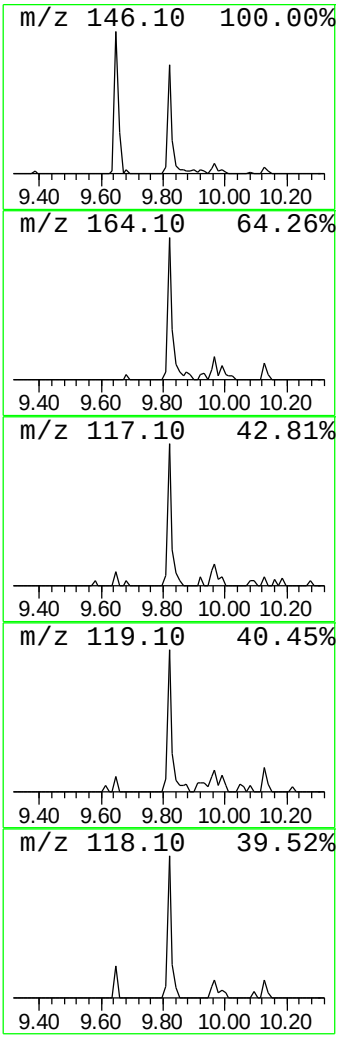
Peak Number 16 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	7.51 ng	128437	Acenaphthene-d10	10.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	95
2	Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	49
3	2-Ethyl-2-phenylaziridine	147	C10H13N	000768-82-1	45
4	2-Hydroxy-1,7-naphthyridine	146	C8H6N2O	054920-82-0	27
5	1,3-Oxazolidine, 4-methyl-2-(4-m...	253	C17H19NO	1000138-83-3	27





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Instrument : BNA_F
ClientSampleId : MW-1

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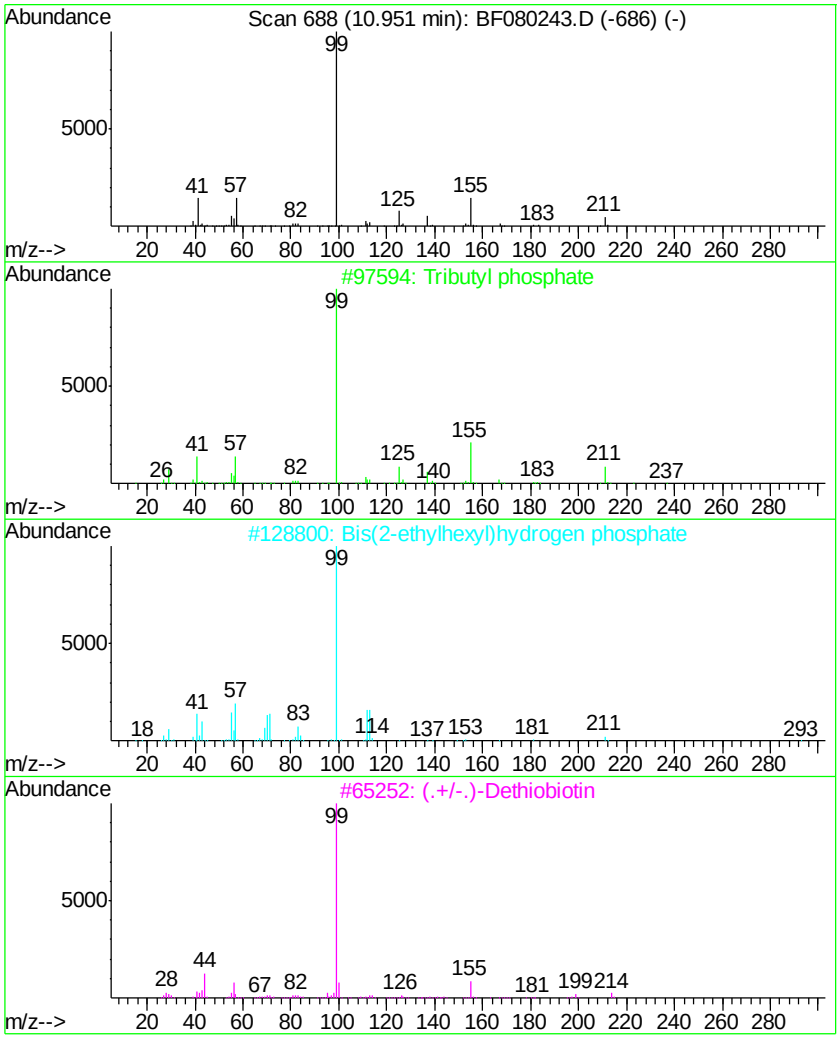
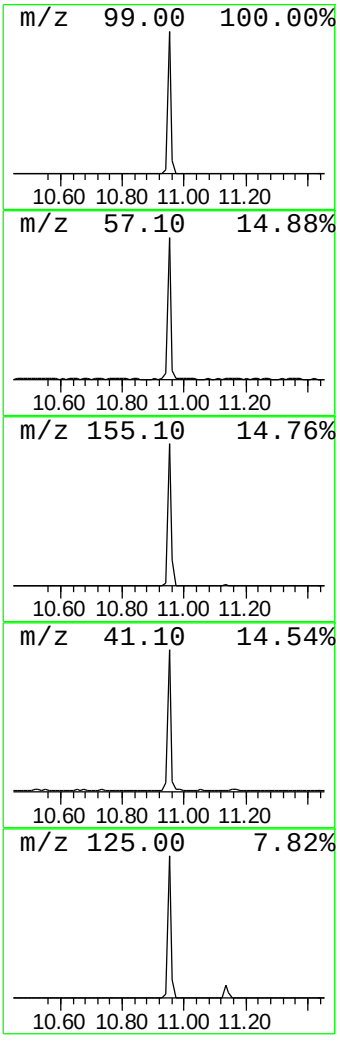
Peak Number 17 Tributyl phosphate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.95	28.90 ng	494284	Acenaphthene-d10	10.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tributyl phosphate	266	C12H27O4P	000126-73-8	91
2	Bis(2-ethylhexyl)hydrogen phosphate	322	C16H35O4P	000298-07-7	36
3	(.+/-.)-Dethiobiotin	214	C10H18N2O3	000636-20-4	36
4	Phosphoric acid, tris(2-ethylhex...	434	C24H51O4P	000078-42-2	36
5	n-Pentyl methylphosphonofluoridate	168	C6H14FO2P	013454-59-6	9





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Sample : G2837-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-1

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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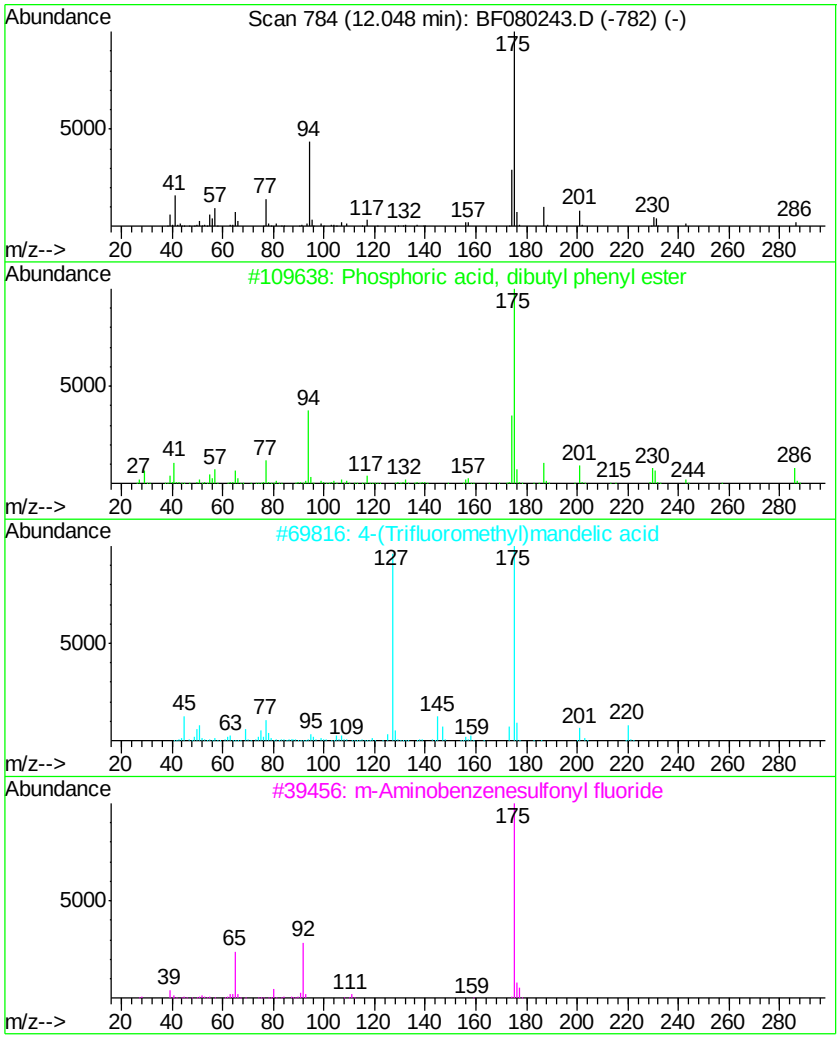
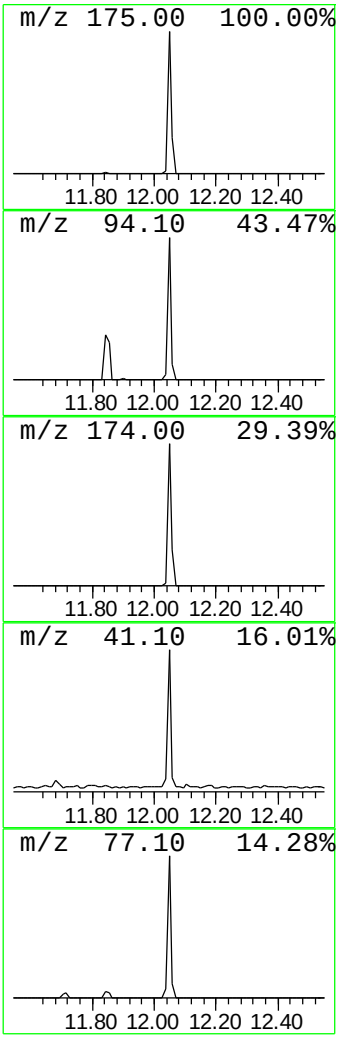
Peak Number 18 Phosphoric acid, dibutyl ph... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	12.99 ng	218996	Phenanthrene-d10	11.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, dibutyl phenyl ...	286	C14H23O4P	002528-36-1	97
2	4-(Trifluoromethyl)mandelic acid	220	C9H7F3O3	000395-35-7	32
3	m-Aminobenzenesulfonyl fluoride	175	C6H6FNO2S	000368-50-3	17
4	2,3,4,5-Tetrafluorobenzonitrile	175	C7HF4N	016582-93-7	9
5	7-Methoxy-2,3-dimethylindole	175	C11H13NO	053918-89-1	9





Tentatively Identified Compound (LSC) summary

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Misc :
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Instrument : BNA_F
ClientSampleId : MW-1

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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
2-Butanol, 2,3-di...	3.74	12.0	ng	126434	1	7.28	210629	20.0
2-Pentanol, 2,4-d...	4.82	7.9	ng	82932	1	7.28	210629	20.0
2-Pentanone, 4-hy...	5.50	4.5	ng	47278	1	7.28	210629	20.0
Benzene, 1-ethyl-...	6.97	4.1	ng	43620	1	7.28	210629	20.0
unknown7.05	7.05	84.3	ng	888167	1	7.28	210629	20.0
unknown7.34	7.34	4.8	ng	50128	1	7.28	210629	20.0
unknown7.52	7.52	11.7	ng	123377	1	7.28	210629	20.0
unknown8.13	8.13	5.3	ng	76678	2	8.57	290056	20.0
Benzene, 2-etheny...	8.31	2.4	ng	34030	2	8.57	290056	20.0
1-Adamantanol	9.11	2.5	ng	36668	2	8.57	290056	20.0
unknown9.52	9.52	2.6	ng	45263	3	10.34	342107	20.0
Benzoic acid, 2,4...	9.82	7.5	ng	128437	3	10.34	342107	20.0
Tributyl phosphate	10.95	28.9	ng	494284	3	10.34	342107	20.0
Phosphoric acid, ...	12.05	13.0	ng	218996	4	11.85	337065	20.0