LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

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-BF110715.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	4.967	248	252	255	rBV	792958	924221	16.76%	2.154%
2	5.424	289	292	294	rBV	2086469	2510356	45.51%	5.849%
3	6.453	380	382	385	rBV	1998588	1694834	30.73%	3.949%
4	6.579	390	393	395	rBV	4781622	4196143	76.08%	9.778%
5	6.807	410	413	415	rBV	962344	1187208	21.53%	2.766%
6	6.876	417	419	424	rBV	386677	430863	7.81%	1.004%
7	6.956	424	426	429	rVB	2697983	3710294	67.27%	8.645%
8	7.367	459	462	464	rBV	3029198	3052375	55.34%	7.112%
9	7.836	499	503	507	rBV2	55497	106594	1.93%	0.248%
10	8.008	515	518	519	rBV	62859	70540	1.28%	0.164%
11	8.088	523	525	528	rBV	1373631	1471996	26.69%	3.430%
12	8.476	556	559	560	rVB3	48300	77598	1.41%	0.181%
13	8.533	560	564	565	rBV3	37458	77345	1.40%	0.180%
14	8.568	565	567	570	rVB2	99422	118438	2.15%	0.276%
15	8.670	573	576	579	rVB4	37335	77174	1.40%	0.180%
16	8.728	579	581	584	rBV3	38784	70656	1.28%	0.165%
17	8.876	587	594	596	rBV4	84146	263870	4.78%	0.615%
18	8.956	599	601	604	rVB3	90080	129958	2.36%	0.303%
19	9.048	604	609	611	rBV5	126926	288937	5.24%	0.673%
20	9.093	611	613	614	rBV	89178	117064	2.12%	0.273%
21	9.173	617	620	622	rBV	6199581	5515474	100.00%	12.852%
22	9.276	625	629	633	rBV6	95712	275106	4.99%	0.641%
23	9.356	633	636	637	rBV2	93464	140480	2.55%	0.327%
24	9.425	640	642	643	rBV	116612	126475	2.29%	0.295%
25	9.493	646	648	651	rBV3	79082	211244	3.83%	0.492%
26	9.562	652	654	657	rVB2	133261	233051	4.23%	0.543%
27	9.665	661	663	665	rBV3	79621	135148	2.45%	0.315%
28	9.722	665	668	670	rBV3	51570	111703	2.03%	0.260%
29	9.791	670	674	676	rBV2	175363	252542	4.58%	0.588%
30	9.848	676	679	681	rBV	1823535	1645954	29.84%	3.835%
31	9.916	682	685	686	rBV	287626	408340	7.40%	0.951%
32	10.088	699	700	701	rBV	95580	68961	1.25%	0.161%
33	10.122	701	703	704	rBV2	94832	106557	1.93%	0.248%
34	10.156	704	706	708	rBV2	66529	92082	1.67%	0.215%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

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-BF110715.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.271	714	716	717	rBV2	98512	118361	2.15%	0.276%
36	10.453	730	732	733	rBV	83772	93494	1.70%	0.218%
37	10.545	738	740	743	rBV3	67598	149492	2.71%	0.348%
38	10.591	743	744	745	rVV	81630	61322	1.11%	0.143%
39	10.636	745	748	750	rVV2	2856493	3209504	58.19%	7.479%
40	10.671	750	751	754	rVB3	100086	129913	2.36%	0.303%
41	10.762	757	759	763	rBV3	174143	283722	5.14%	0.661%
42	11.334	806	809	811	rBV	1402321	1415218	25.66%	3.298%
43	11.882	855	857	860	rVB2	610906	630214	11.43%	1.468%
44	12.008	866	868	870	rBV	183122	181286	3.29%	0.422%
45	12.671	924	926	928	rBV	777638	758709	13.76%	1.768%
46	12.922	945	948	950	rVB	4202408	3860580	70.00%	8.996%
47	13.974	1037	1040	1042	rBV	852601	1098498	19.92%	2.560%
48	15.391	1161	1164	1167	rVB	848847	1026395	18.61%	2.392%

Sum of corrected areas: 42916289

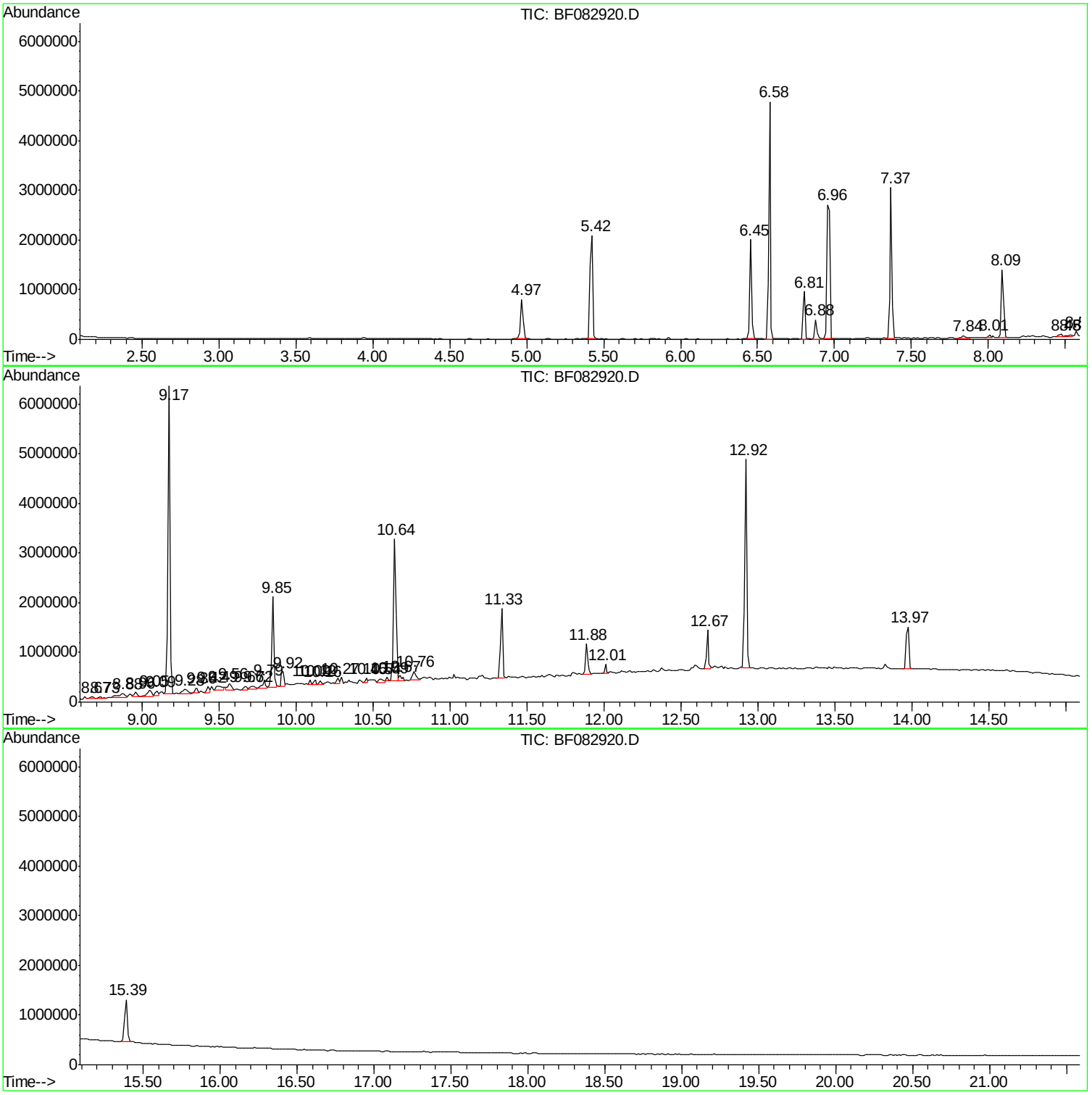
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

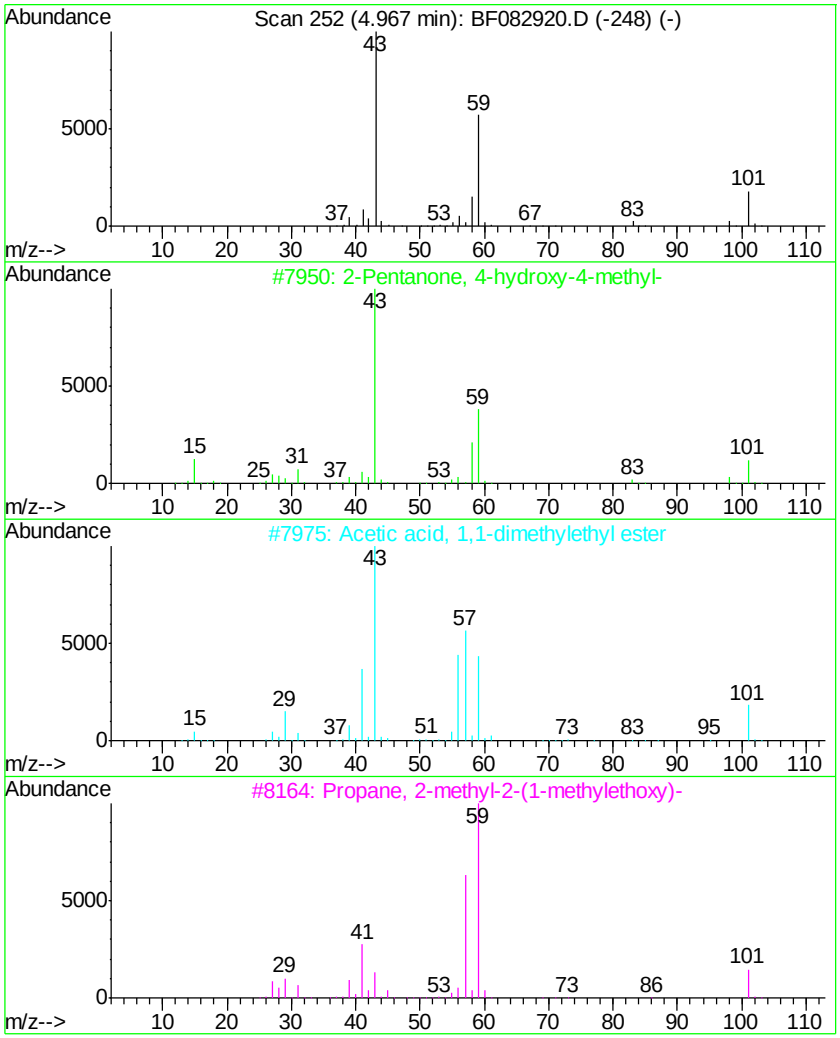
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

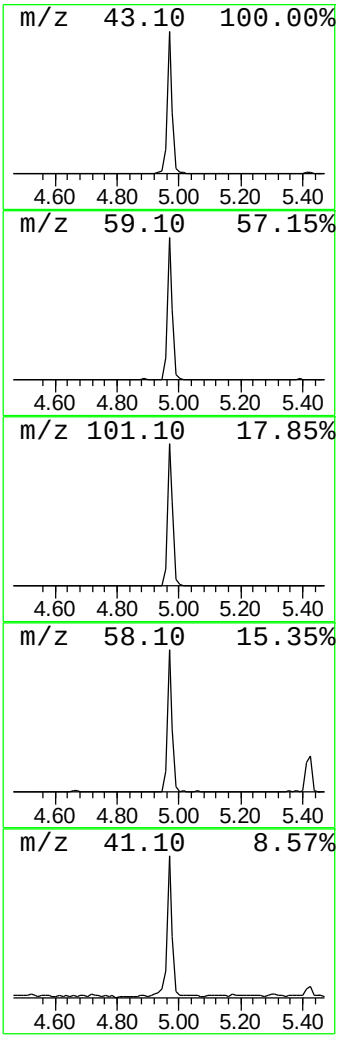
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	15.57 ng	924221	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23





Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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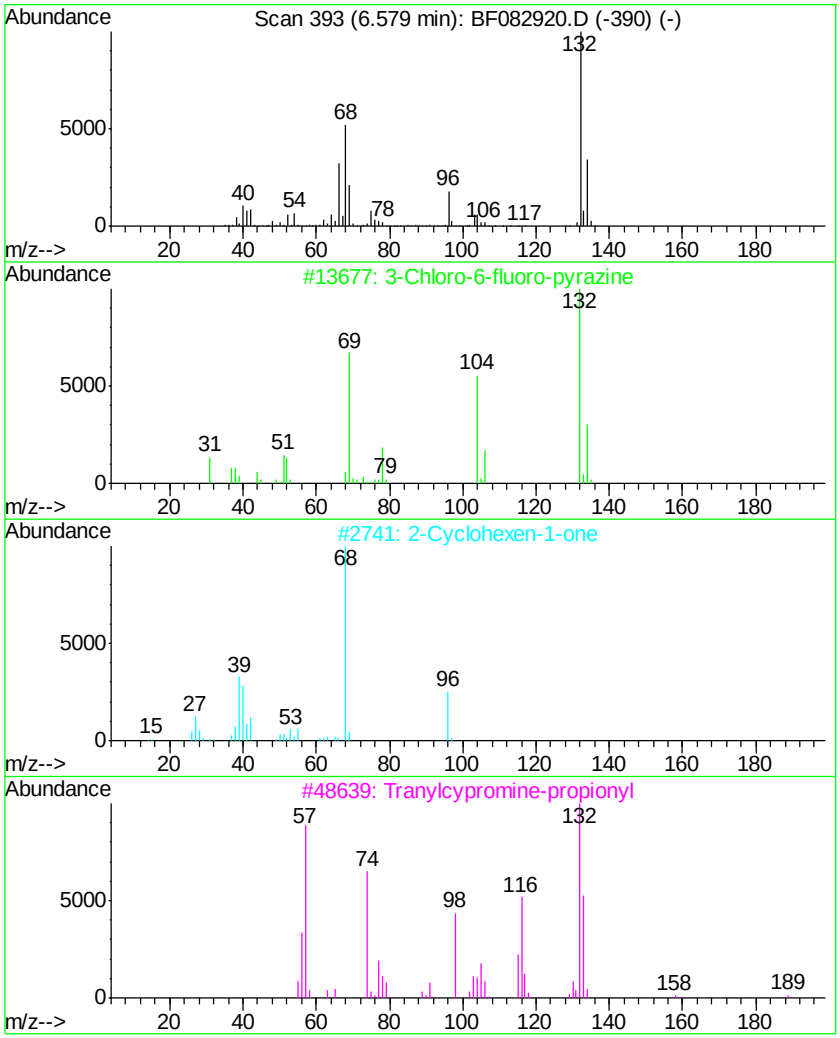
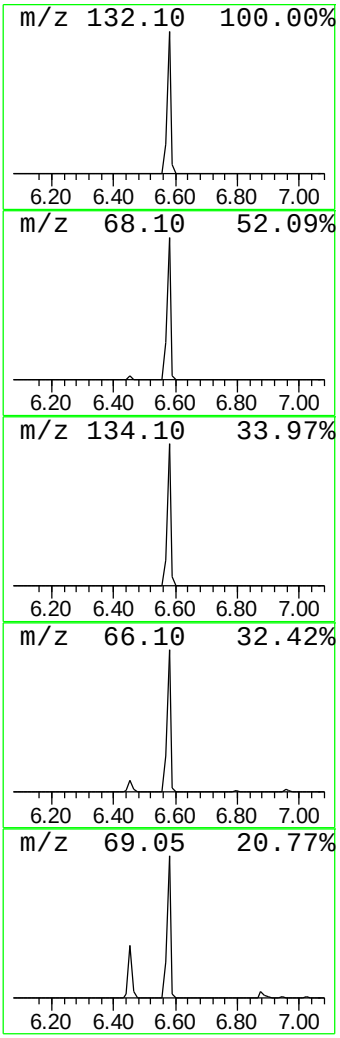
Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	70.69 ng	4196140	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4	Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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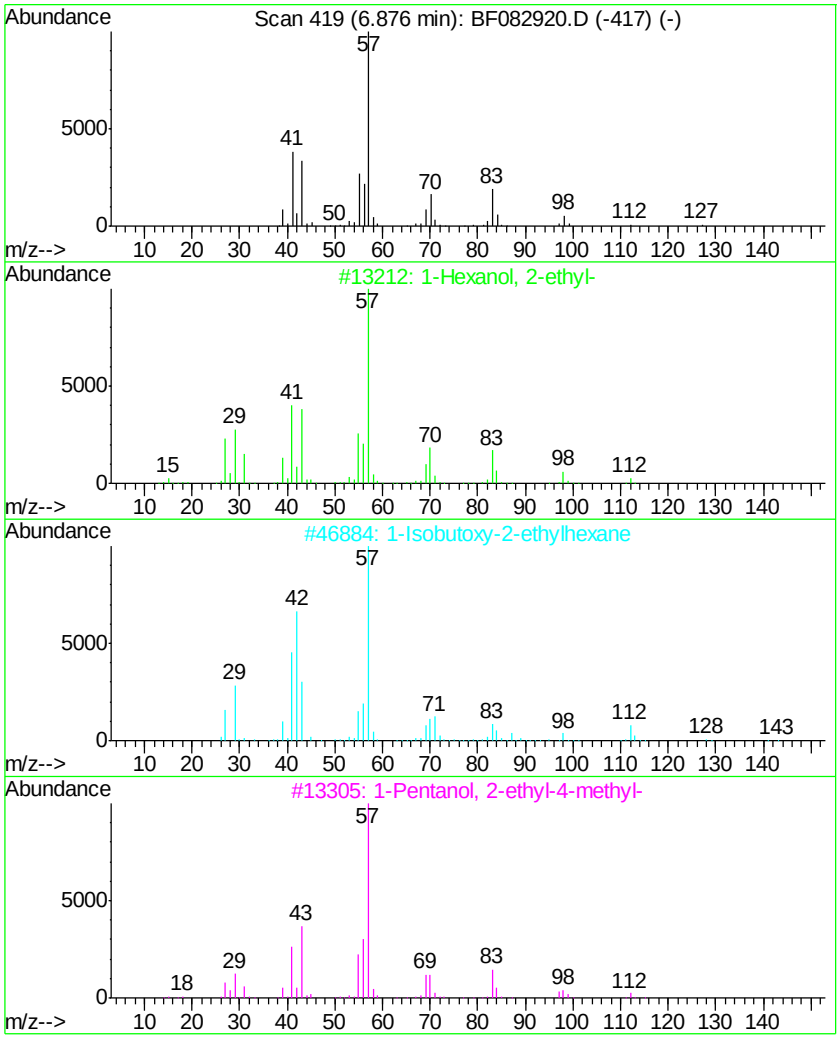
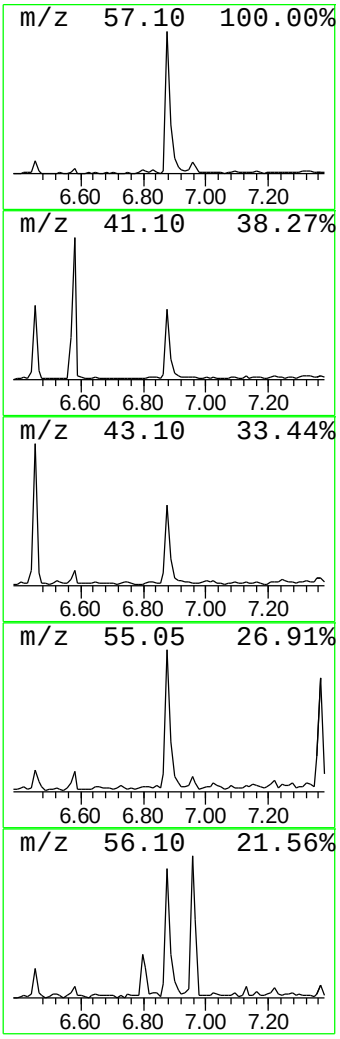
Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	7.26 ng	430863	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	64
3	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	50
4	Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	47
5	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	47





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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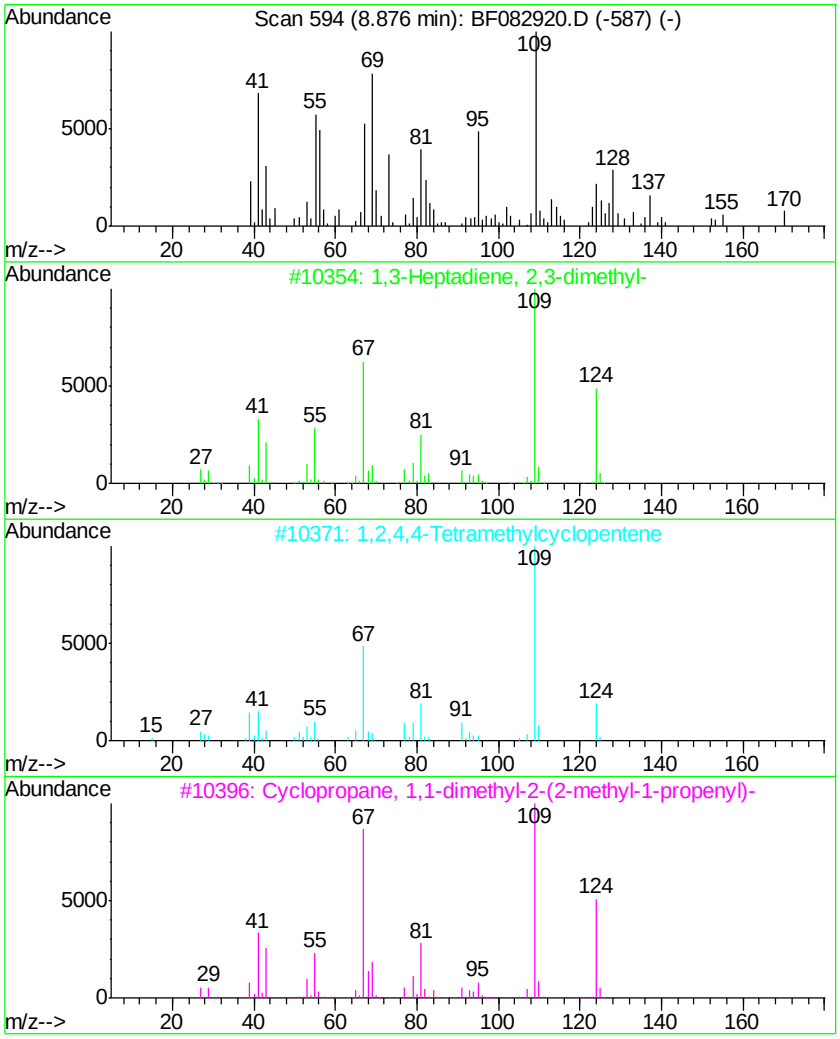
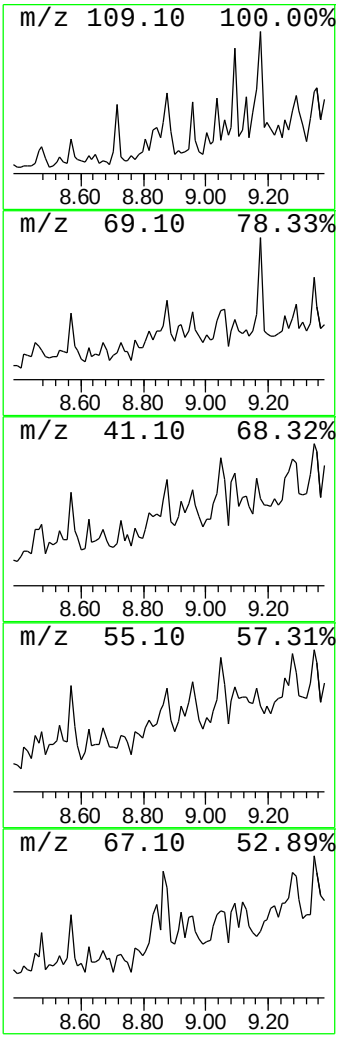
Peak Number 5 unknown8.88 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.88	3.59 ng	263870	Naphthalene-d8	8.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Heptadiene, 2,3-dimethyl-	124	C9H16	074779-65-0	30
2	1,2,4,4-Tetramethylcyclopentene	124	C9H16	065378-76-9	30
3	Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	033422-32-1	27
4	1,1-Dimethyl-4-methylenecyclohexane	124	C9H16	006007-96-1	27
5	1,3-Hexadiene, 2,3,5-trimethyl-	124	C9H16	061142-34-5	27





Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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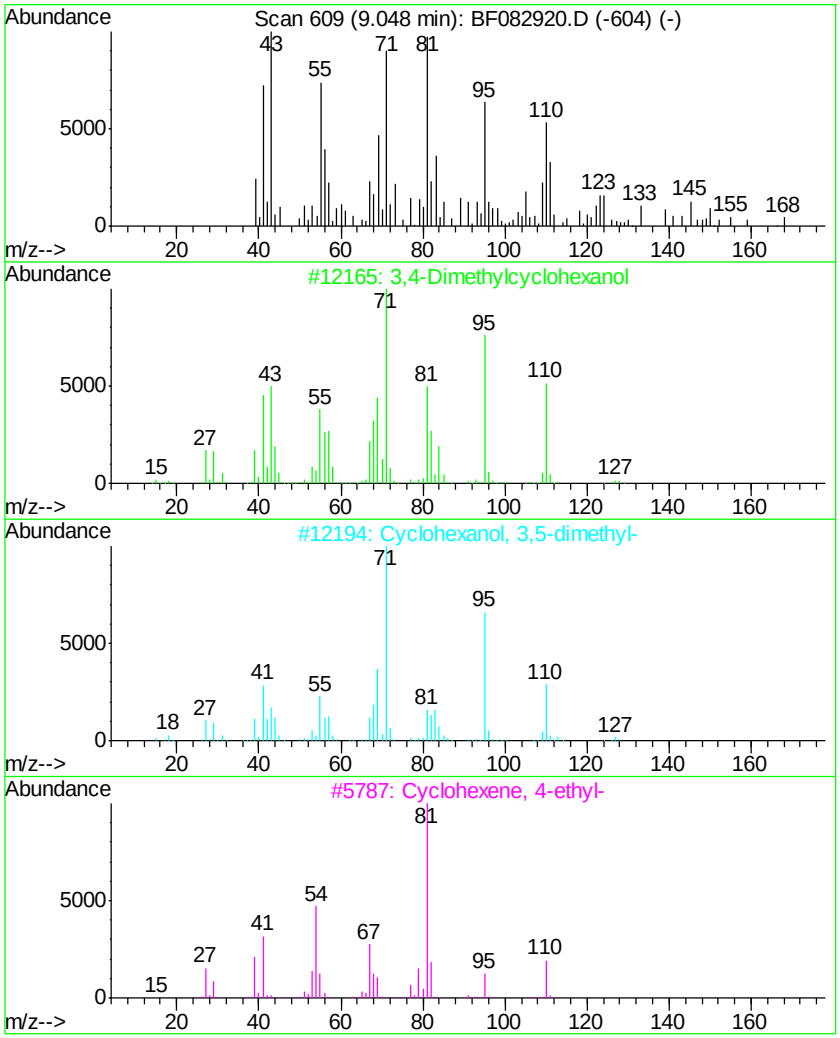
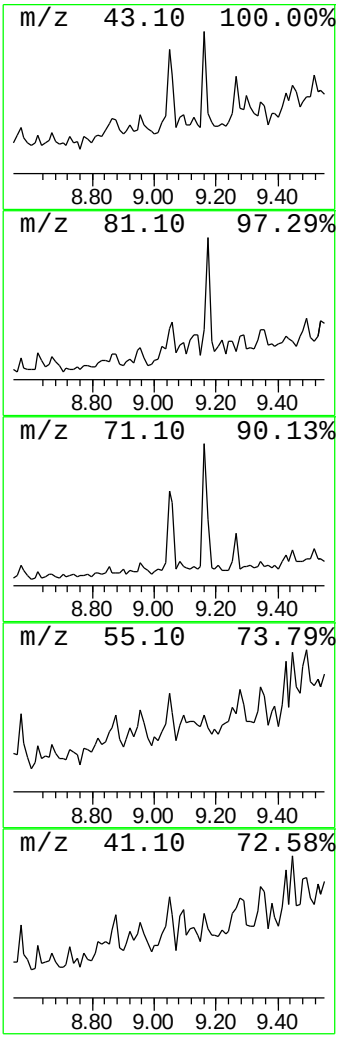
Peak Number 6 unknown9.05 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	3.51 ng	288937	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dimethylcyclohexanol	128	C8H16O	005715-23-1	46
2	Cyclohexanol, 3,5-dimethyl-	128	C8H16O	005441-52-1	43
3	Cyclohexene, 4-ethyl-	110	C8H14	003742-42-5	30
4	2,4-Heptadienal, (E,E)-	110	C7H10O	004313-03-5	25
5	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	22





Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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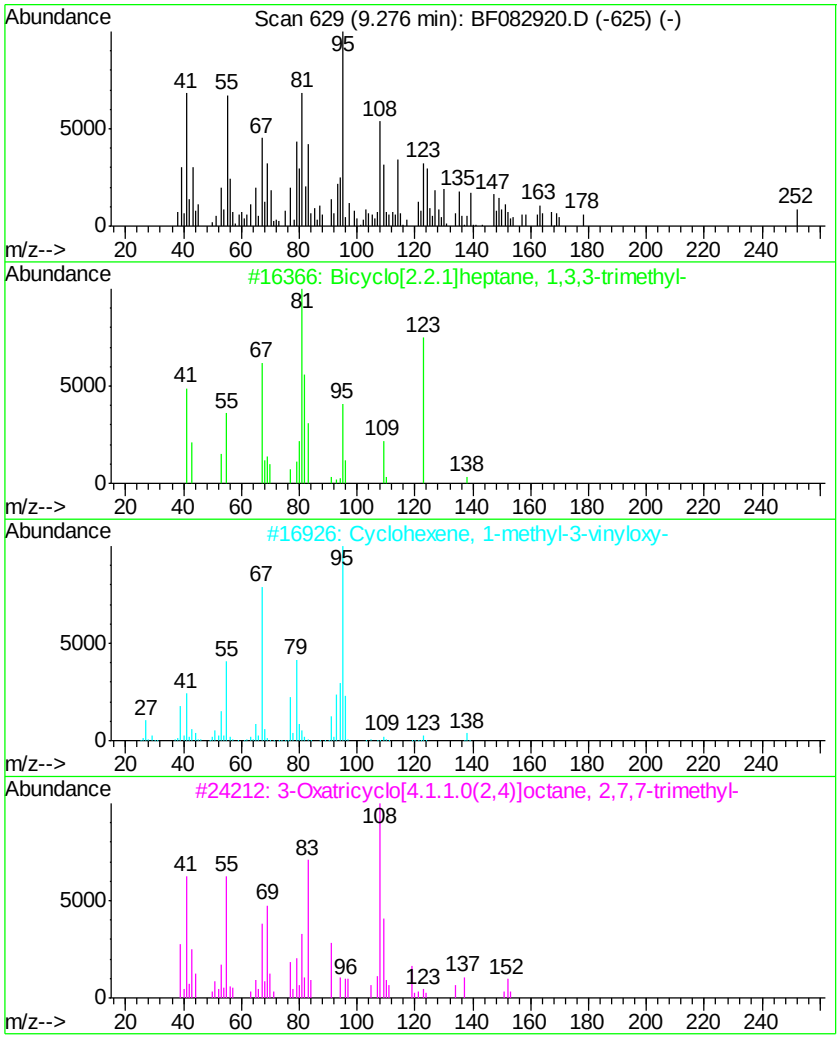
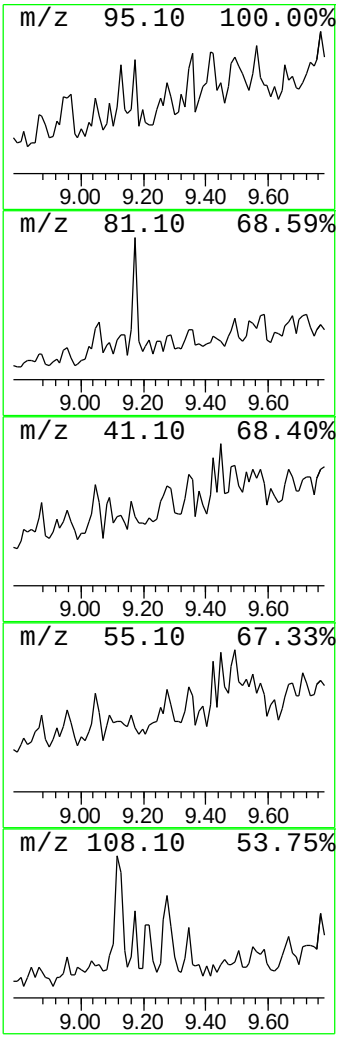
Peak Number 7 unknown9.28 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	3.34 ng	275106	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	27
2	Cyclohexene, 1-methyl-3-vinyloxy-	138	C9H14O	100144-30-7	25
3	3-Oxatricyclo[4.1.1.0(2,4)]octan...	152	C10H16O	001686-14-2	25
4	Cyclohexene, 1-methyl-3-(formylm...	138	C9H14O	129993-40-4	22
5	Cyclohexene, 1-chloro-6-methyl-	130	C7H11Cl	016642-50-5	22





Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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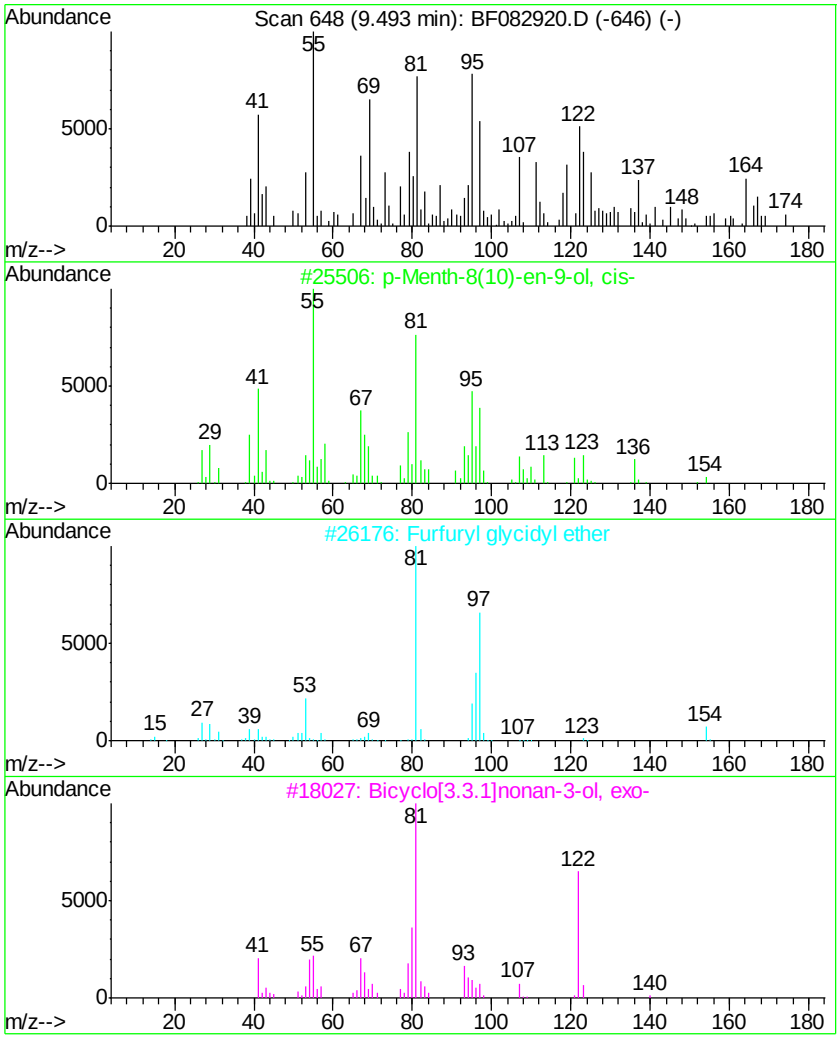
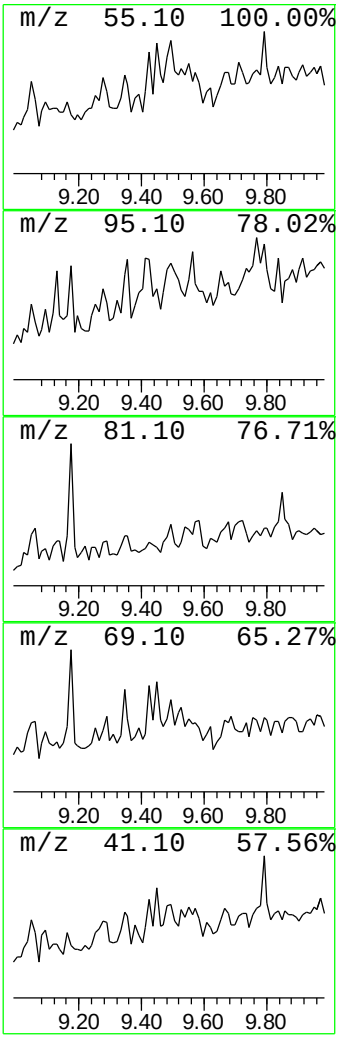
Peak Number 8 unknown9.49 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	2.57 ng	211244	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Menth-8(10)-en-9-ol, cis-	154	C10H18O	015714-13-3	35
2	Furfuryl glycidyl ether	154	C8H10O3	005380-87-0	25
3	Bicyclo[3.3.1]nonan-3-ol, exo-	140	C9H16O	010036-08-5	22
4	Bicyclo[3.3.1]nonan-9-ol	140	C9H16O	015598-80-8	18
5	5-Hepten-2-ol, 6-methyl-	128	C8H16O	001569-60-4	11





Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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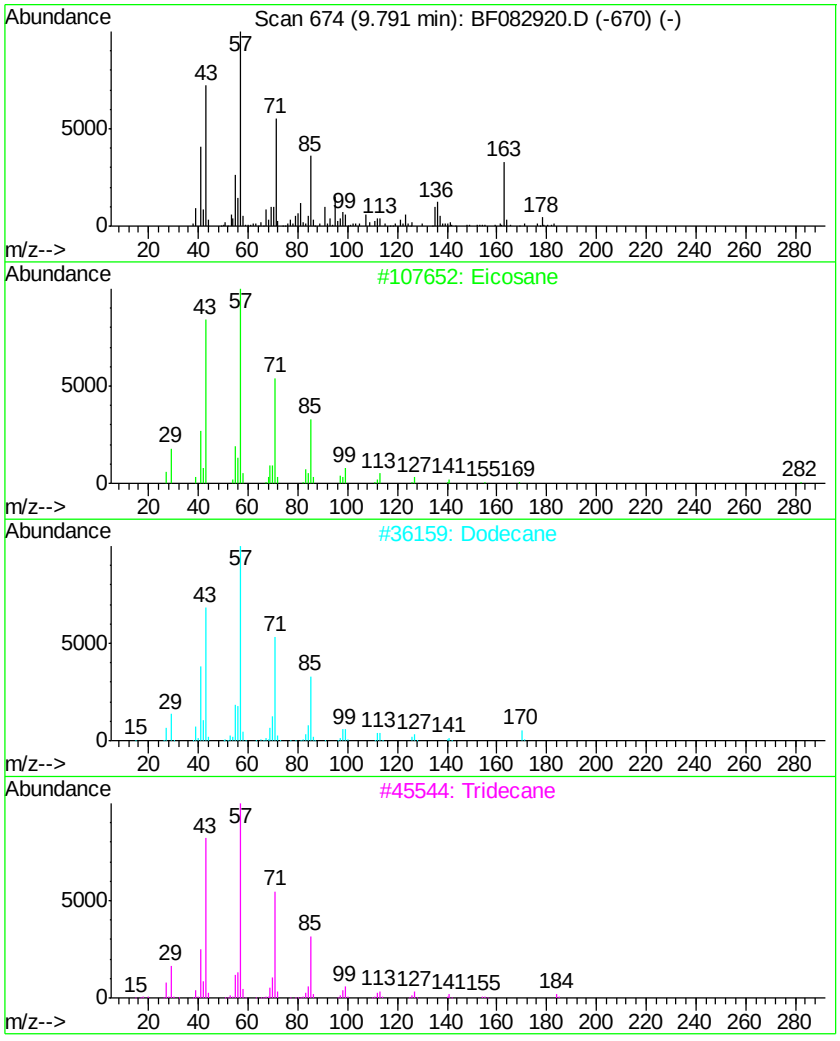
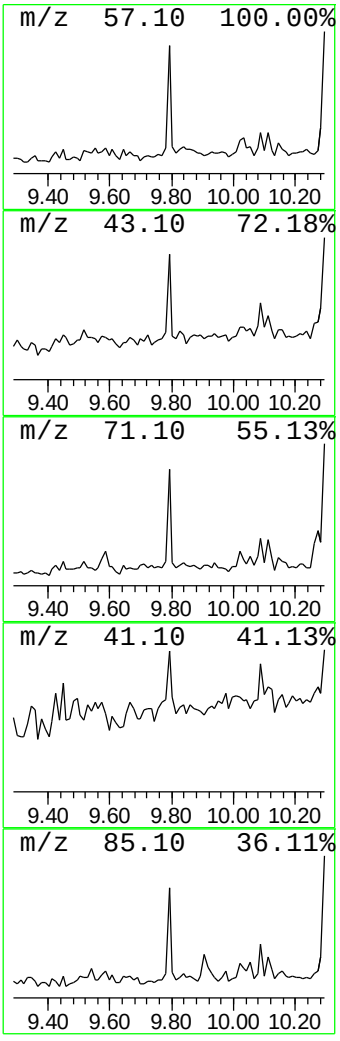
Peak Number 10 unknown9.79 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	3.07 ng	252542	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	49
2	Dodecane	170	C12H26	000112-40-3	46
3	Tridecane	184	C13H28	000629-50-5	46
4	Pentadecane	212	C15H32	000629-62-9	43
5	Hexadecane	226	C16H34	000544-76-3	43





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Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.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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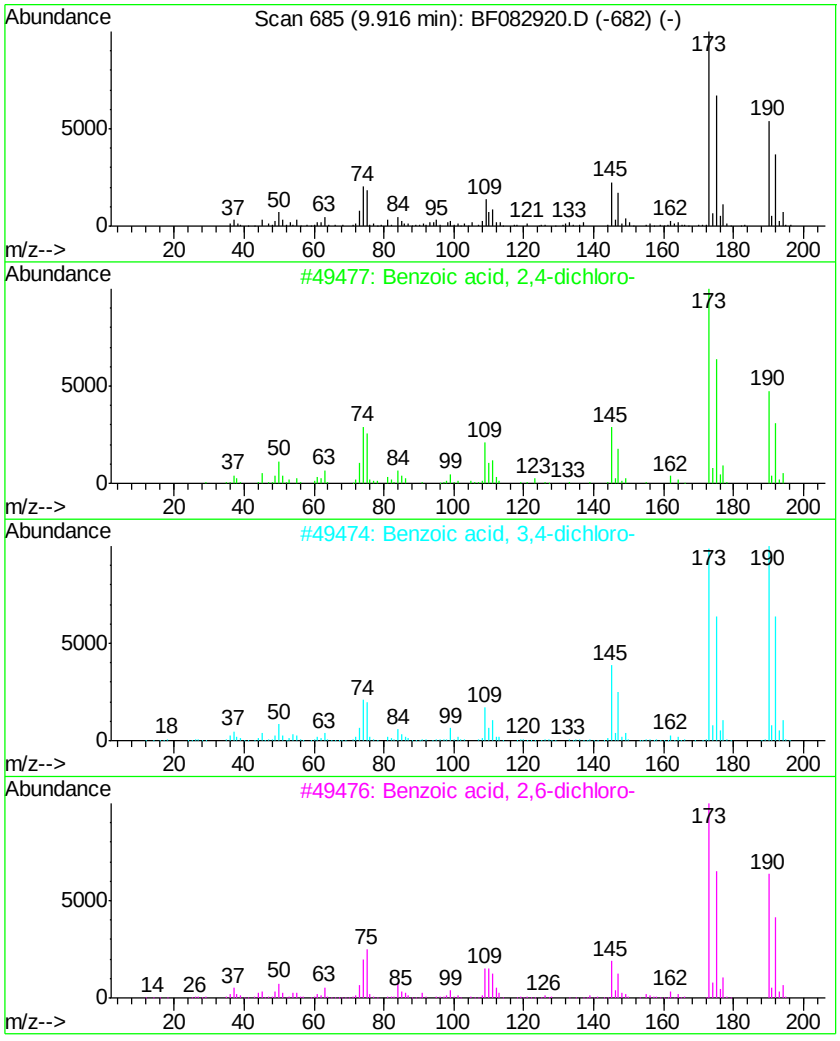
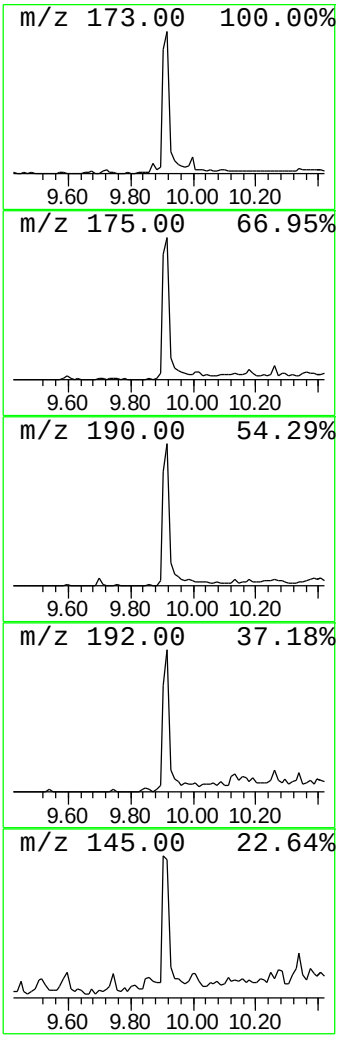
Peak Number 11 Benzoic acid, 2,4-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.92	4.96 ng	408340	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, 2,4-dichloro-	190	C7H4Cl2O2	000050-84-0	97
2	Benzoic acid, 3,4-dichloro-	190	C7H4Cl2O2	000051-44-5	95
3	Benzoic acid, 2,6-dichloro-	190	C7H4Cl2O2	000050-30-6	94
4	3,5-Dichlorobenzoic acid	190	C7H4Cl2O2	000051-36-5	93
5	Benzoic acid, 2,5-dichloro-	190	C7H4Cl2O2	000050-79-3	92





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111315\
Data File : BF082920.D
Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
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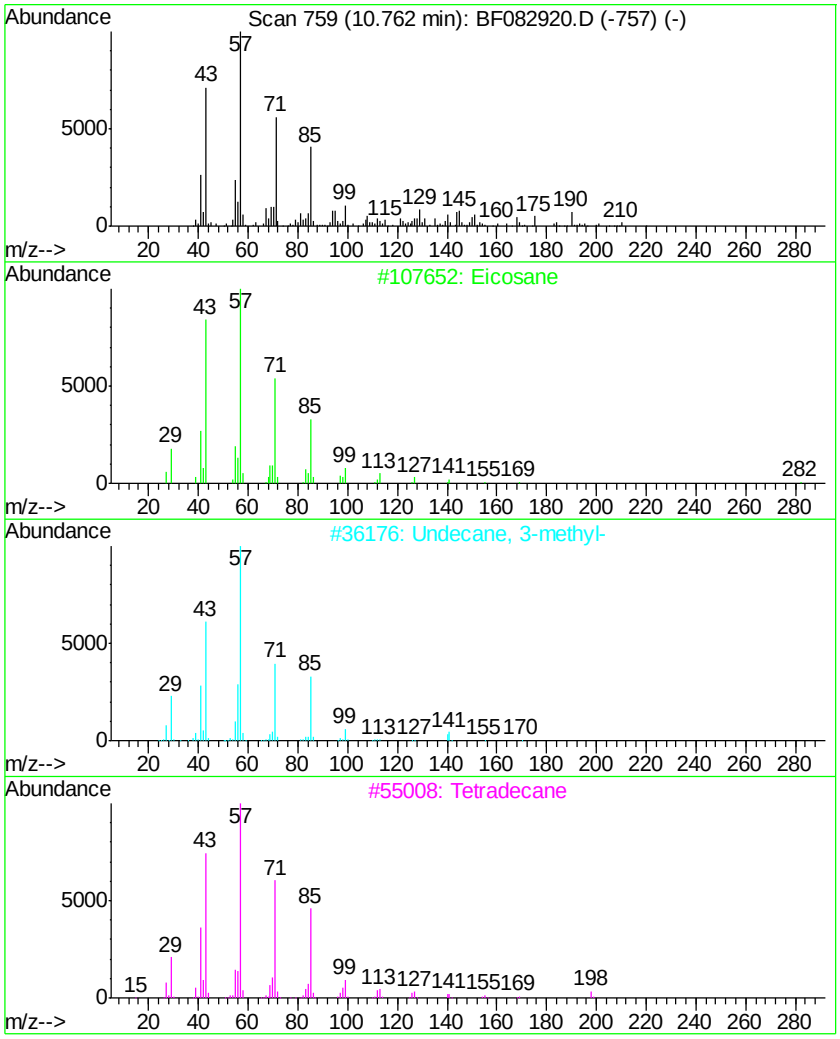
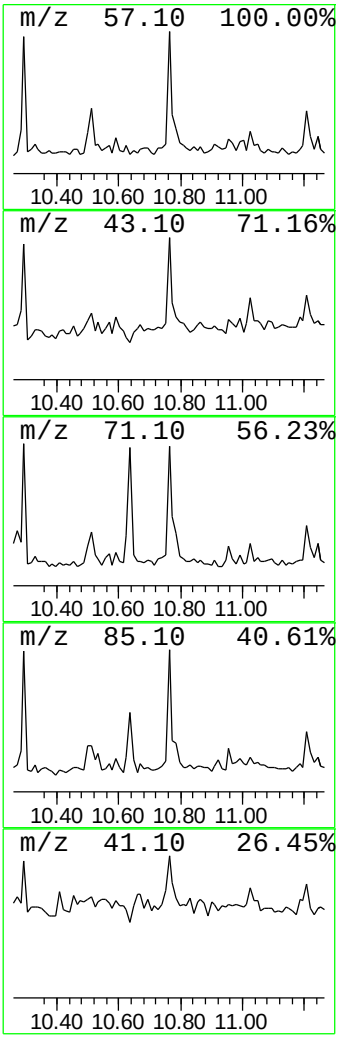
Peak Number 12 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	4.01 ng	283722	Phenanthrene-d10	11.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	74
2	Undecane, 3-methyl-	170	C12H26	001002-43-3	64
3	Tetradecane	198	C14H30	000629-59-4	64
4	Pentadecane	212	C15H32	000629-62-9	64
5	Hexadecane	226	C16H34	000544-76-3	64





Library Search Compound Report

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Acq On : 14 Nov 2015 3:17
Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
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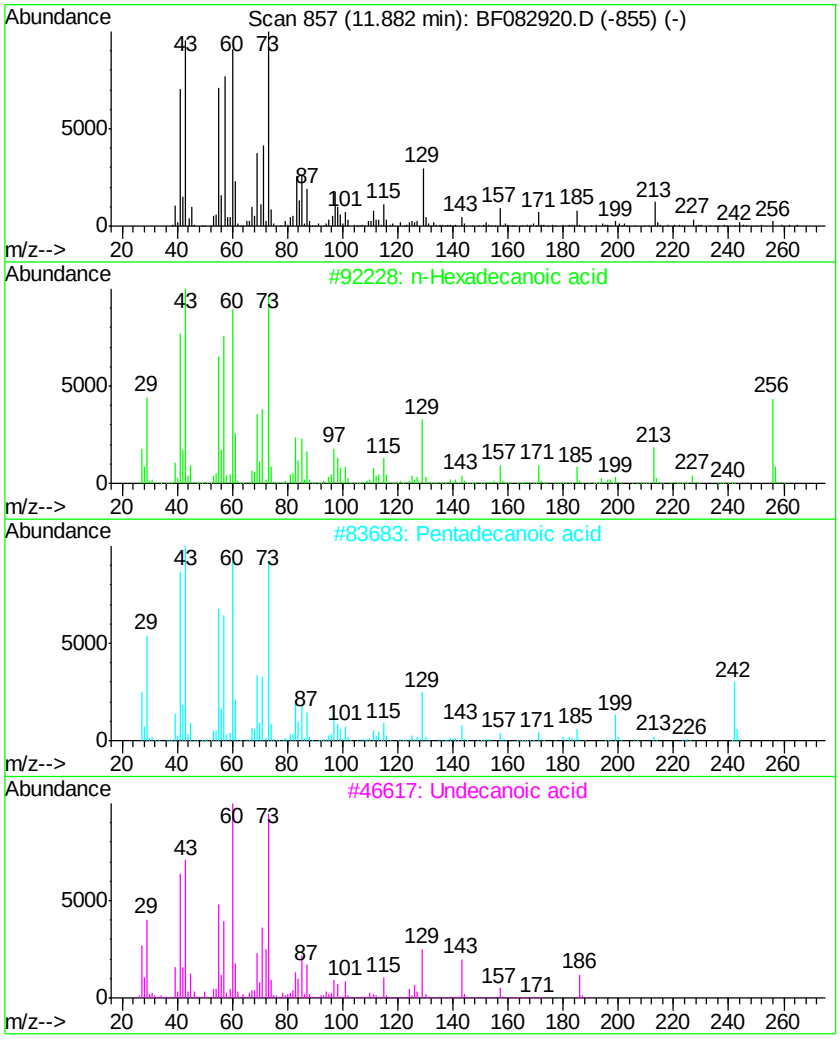
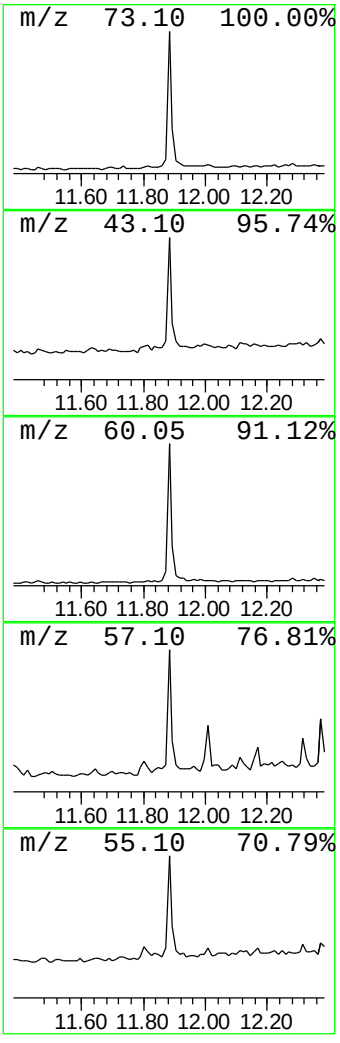
Peak Number 13 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	8.91 ng	630214	Phenanthrene-d10	11.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	Undecanoic acid	186	C11H22O2	000112-37-8	86
4	Dodecanoic acid	200	C12H24O2	000143-07-7	86
5	Tridecanoic acid	214	C13H26O2	000638-53-9	72





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Operator : UM/IZ
Sample : G4382-03
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : URS MW-13

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
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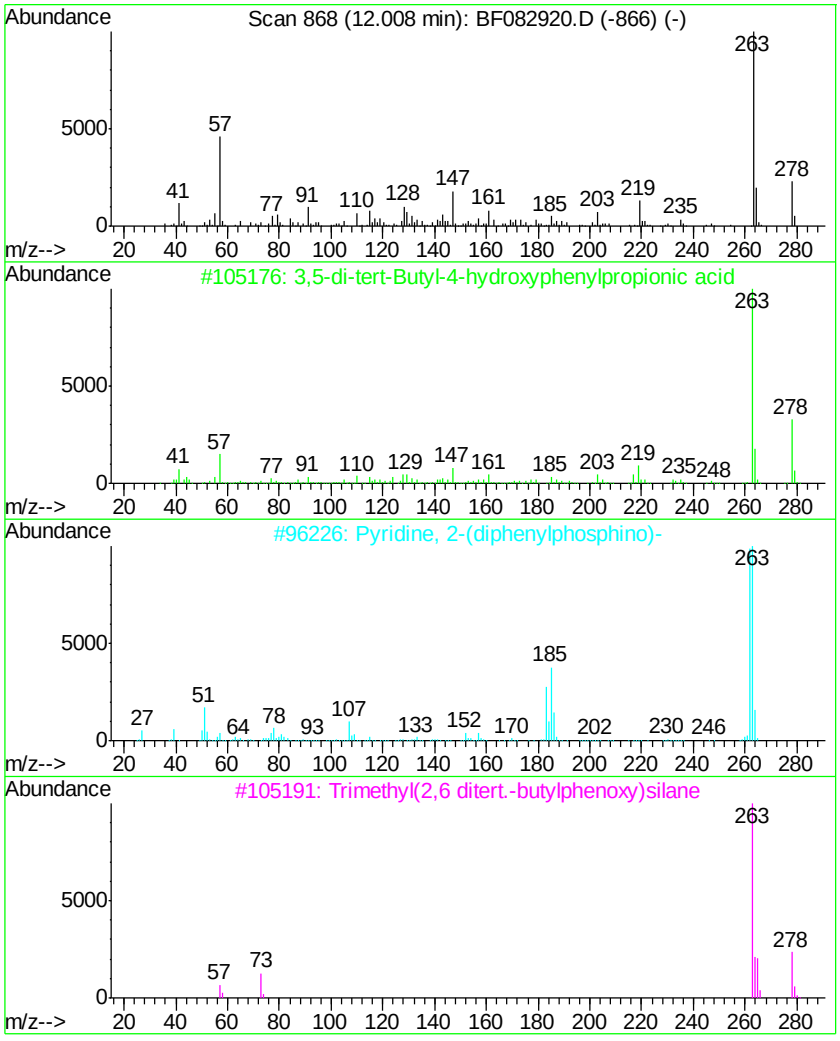
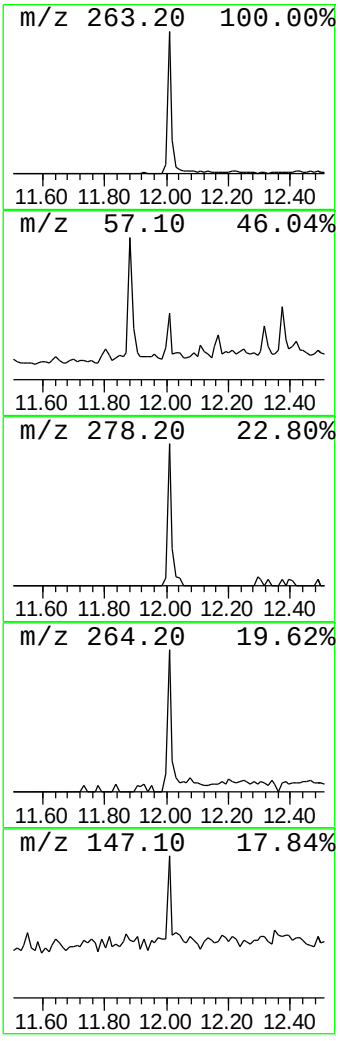
Peak Number 14 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	2.56 ng	181286	Phenanthrene-d10	11.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-di-tert-Butyl-4-hydroxyphenyl...	278	C17H26O3	020170-32-5	94
2	Pyridine, 2-(diphenylphosphino)-	263	C17H14NP	037943-90-1	47
3	Trimethyl(2,6 ditert.-butylpheno...	278	C17H30OSi	010416-73-6	45
4	9(10H)-Acridone, 4,5-dichloro-	263	C13H7Cl2NO	1000163-26-5	43
5	12-Cyclohex-3-enyl-3-methyl-8,9,...	344	C23H24N2O	1000296-08-5	38





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Sample : G4382-03
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ALS Vial : 26 Sample Multiplier: 1

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ClientSampleId : URS MW-13

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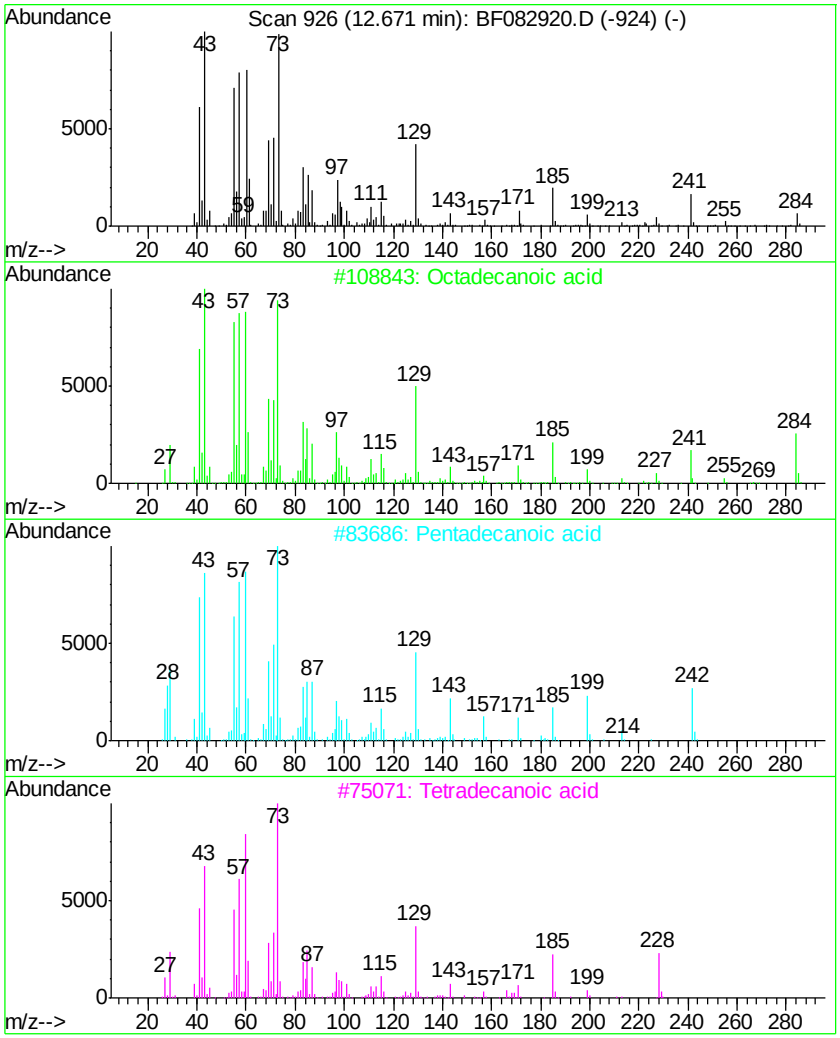
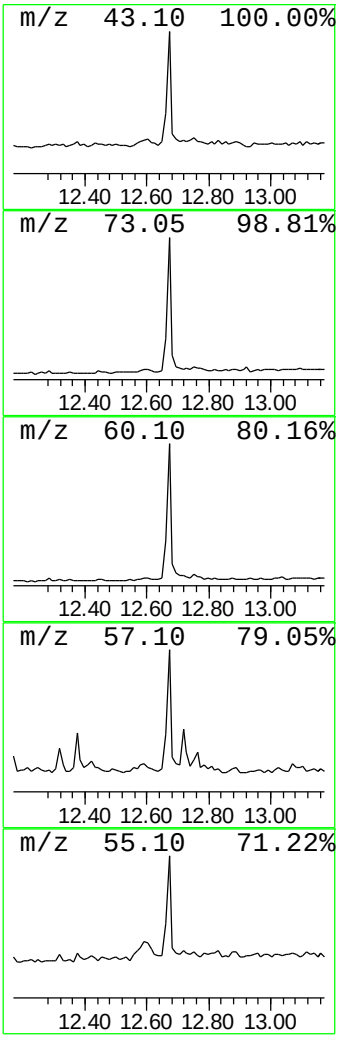
Peak Number 15 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	13.81 ng	758709	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4	Eicosane	282	C20H42	000112-95-8	25
5	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	11





Tentatively Identified Compound (LSC) summary

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					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.97	15.6	ng	924221	1	6.81	1187210	20.0
unknown6.58	6.58	70.7	ng	4196140	1	6.81	1187210	20.0
1-Hexanol, 2-ethyl-	6.88	7.3	ng	430863	1	6.81	1187210	20.0
unknown8.88	8.88	3.6	ng	263870	2	8.09	1472000	20.0
unknown9.05	9.05	3.5	ng	288937	3	9.85	1645950	20.0
unknown9.28	9.28	3.3	ng	275106	3	9.85	1645950	20.0
unknown9.49	9.49	2.6	ng	211244	3	9.85	1645950	20.0
unknown9.79	9.79	3.1	ng	252542	3	9.85	1645950	20.0
Benzoic acid, 2,4...	9.92	5.0	ng	408340	3	9.85	1645950	20.0
Eicosane	10.76	4.0	ng	283722	4	11.33	1415220	20.0
n-Hexadecanoic acid	11.88	8.9	ng	630214	4	11.33	1415220	20.0
3,5-di-tert-Butyl...	12.01	2.6	ng	181286	4	11.33	1415220	20.0
Octadecanoic acid	12.67	13.8	ng	758709	5	13.97	1098500	20.0