LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-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-BF101015.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	4.994	247	250	252	rBV	559074	934341	52.01%	6.226%
2	5.417	285	287	289	rBV	1429089	1379012	76.77%	9.188%
3	6.457	376	378	381	rBV	949050	1363414	75.90%	9.084%
4	6.594	387	390	392	rBV	1452335	1554371	86.53%	10.357%
5	6.640	392	394	397	rVB	108576	110758	6.17%	0.738%
6	6.823	408	410	413	rBV	494486	425660	23.70%	2.836%
7	6.983	421	424	426	rBV	1222188	1270724	70.74%	8.467%
8	7.394	458	460	462	rBV	979027	900212	50.11%	5.998%
9	8.115	520	523	525	rBV	654662	606407	33.76%	4.040%
10	8.537	558	560	564	rBV	76145	131177	7.30%	0.874%
11	8.709	573	575	577	rBV	64335	82845	4.61%	0.552%
12	9.132	611	612	615	rBV	150277	117350	6.53%	0.782%
13	9.189	615	617	620	rBV	1359119	1796317	100.00%	11.969%
14	9.269	622	624	626	rBV	156568	161262	8.98%	1.074%
15	9.589	650	652	653	rBV	348069	296942	16.53%	1.979%
16	9.623	653	655	656	rVB2	109964	109229	6.08%	0.728%
17	9.795	669	670	673	rVB2	87600	111735	6.22%	0.744%
18	9.875	673	677	679	rBV	655306	717663	39.95%	4.782%
19	10.126	697	699	701	rBV3	41795	67860	3.78%	0.452%
20	10.206	703	706	708	rBV	166023	176986	9.85%	1.179%
21	10.298	710	714	716	rVV2	83623	128124	7.13%	0.854%
22	10.515	731	733	735	rBV	67420	102466	5.70%	0.683%
23	10.663	744	746	749	rBV	935587	890793	49.59%	5.935%
24	10.778	754	756	760	rVV	247865	376719	20.97%	2.510%
25	10.972	771	773	774	rBV2	50323	91171	5.08%	0.607%
26	11.098	783	784	786	rBV2	43079	75166	4.18%	0.501%
27	11.223	794	795	796	rBV	102108	105511	5.87%	0.703%
28	11.258	796	798	800	rVB	253587	281091	15.65%	1.873%
29	11.361	805	807	810	rBV	584798	642942	35.79%	4.284%

Sum of corrected areas: 15008248

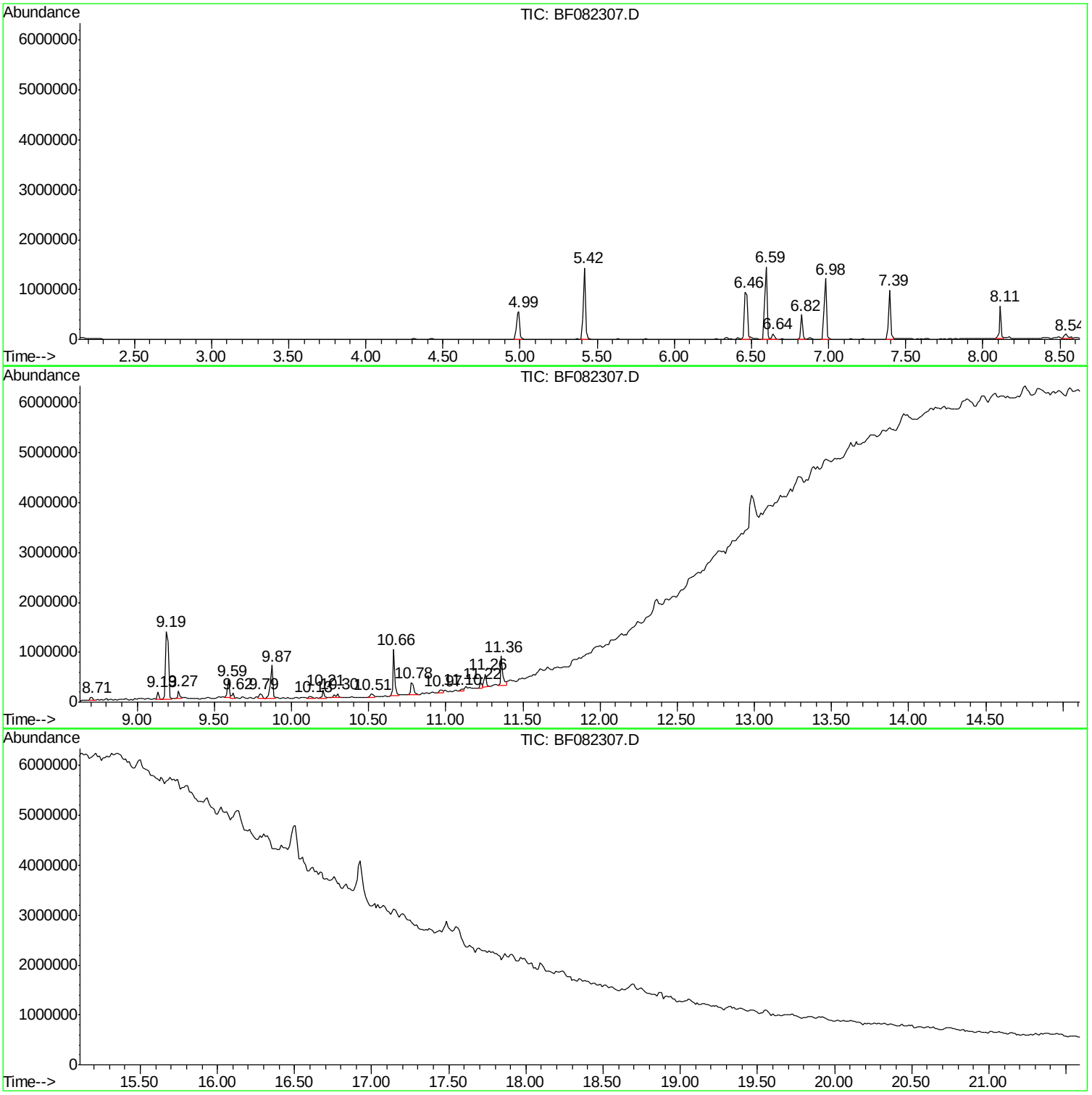
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

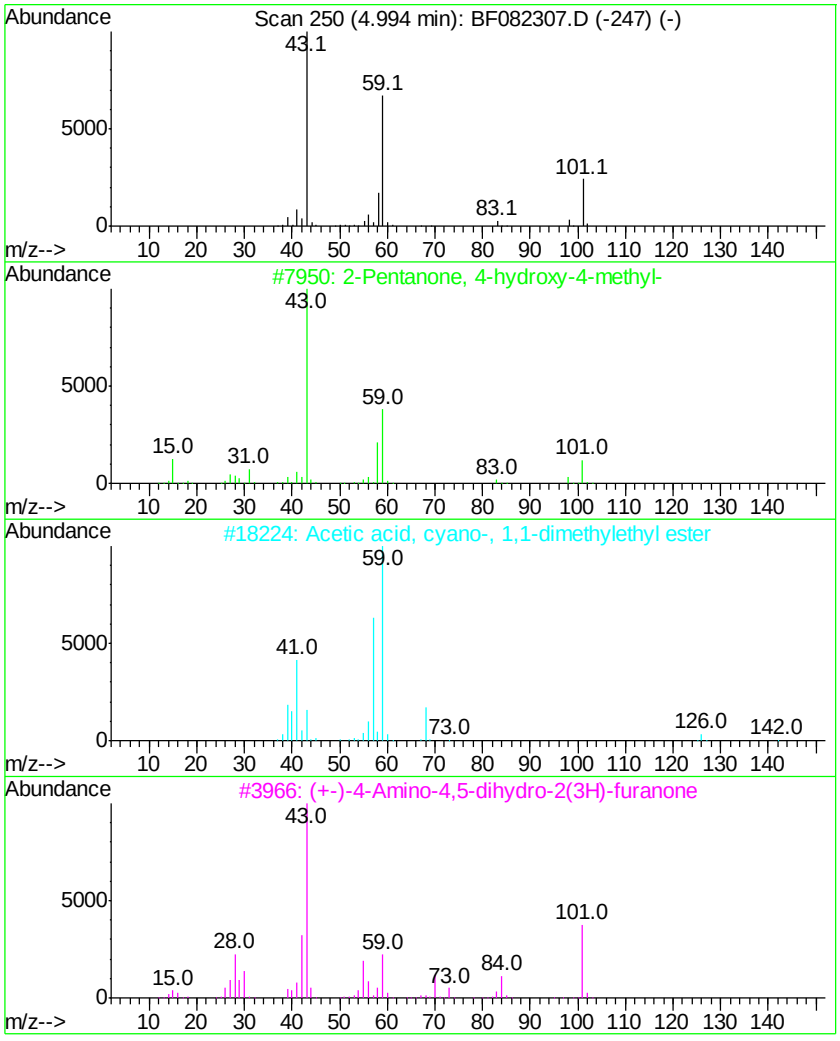
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

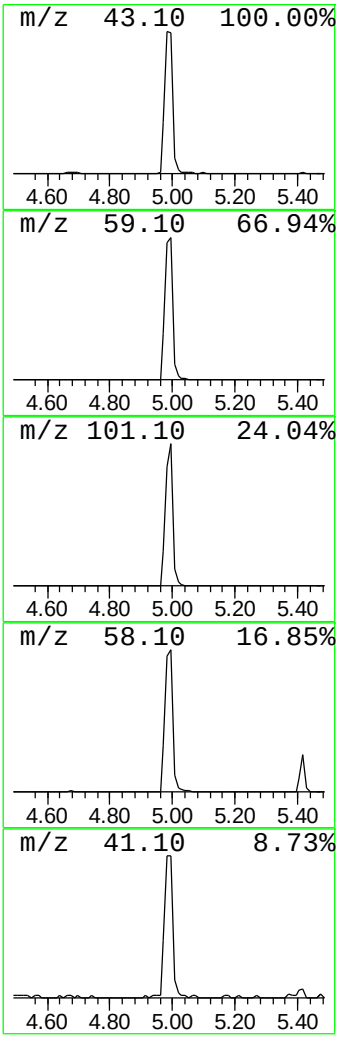
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	43.90 ng	934341	1,4-Dichlorobenzene-d4	6.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Guanidine	59	CH5N3	000113-00-8	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

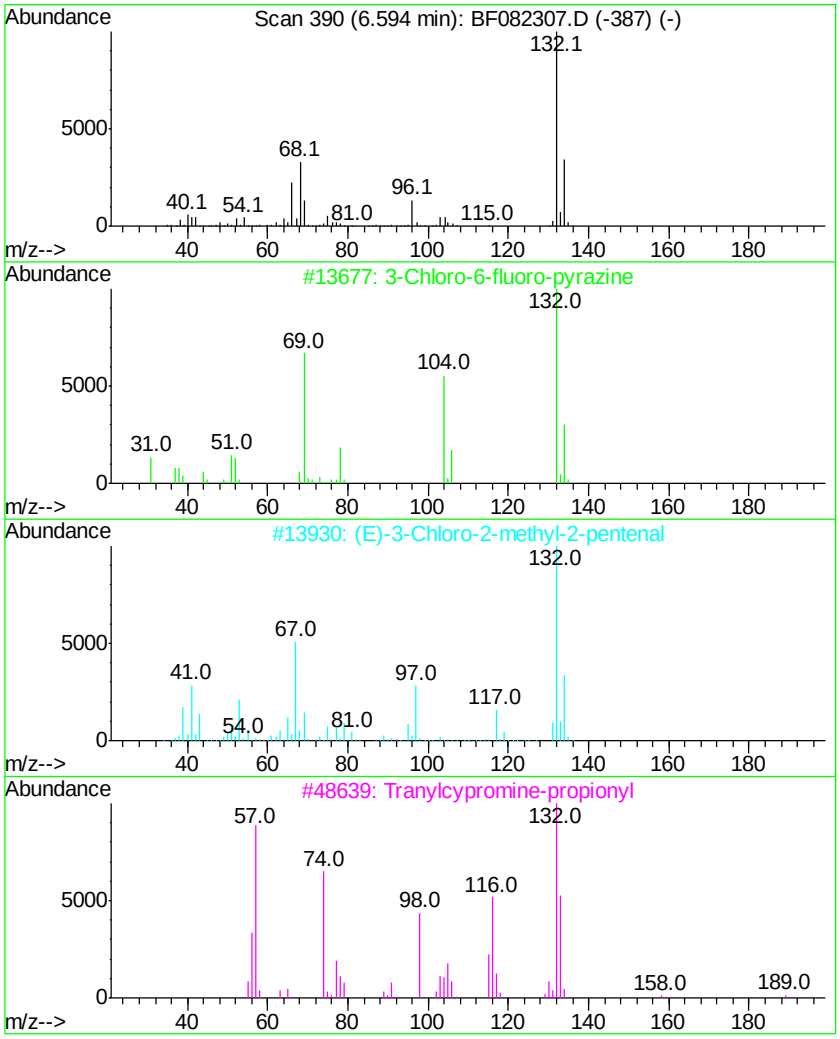
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

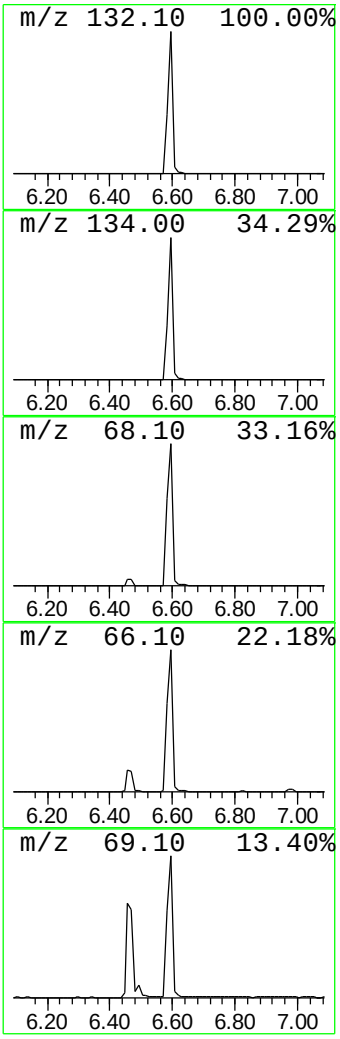
****

Peak Number 2 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	73.03 ng	1554370	1,4-Dichlorobenzene-d4	6.82

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	12
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	5-Aminoindole	132	C8H8N2	005192-03-0	12
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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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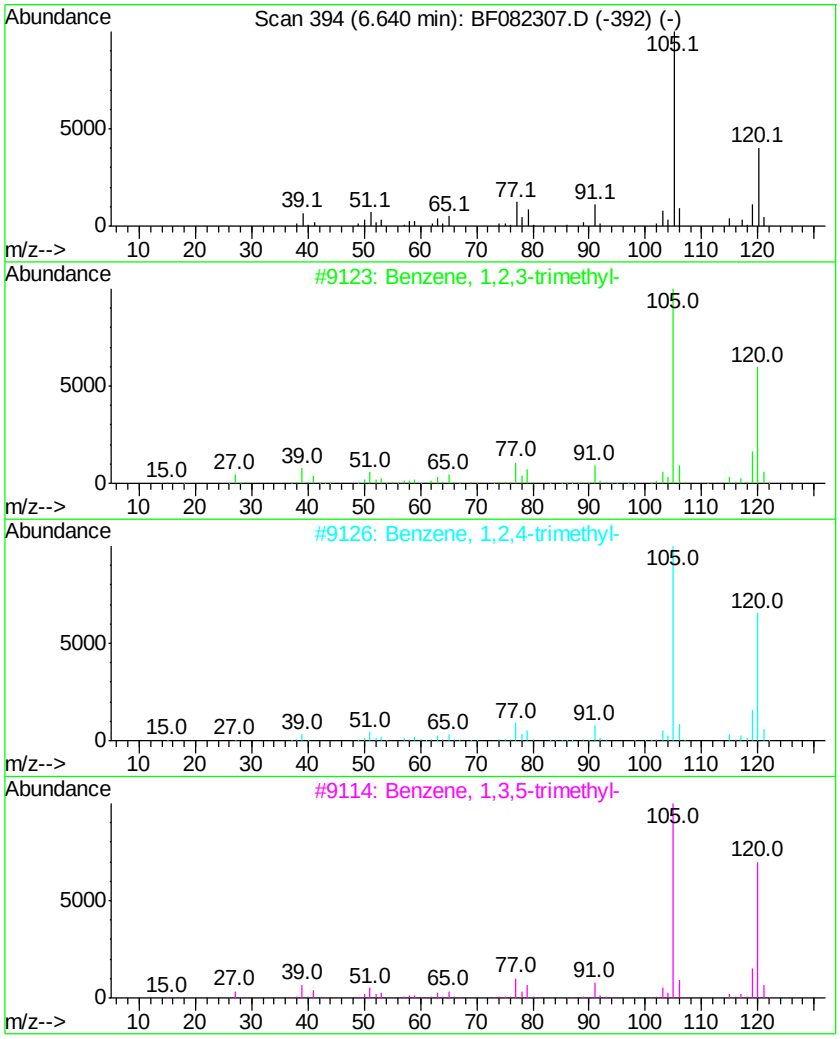
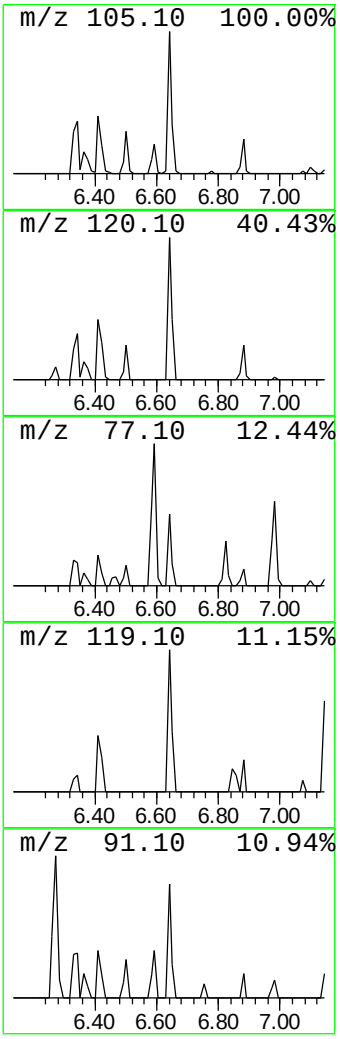
Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	5.20 ng	110758	1,4-Dichlorobenzene-d4	6.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
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-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

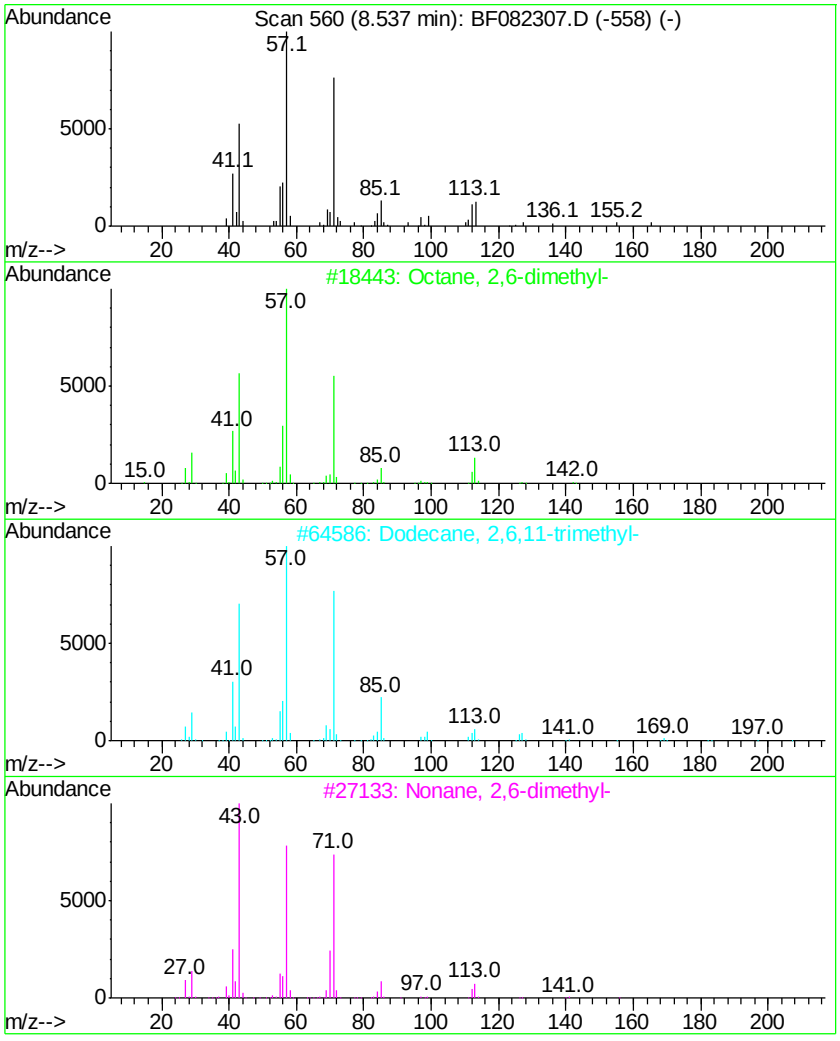
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

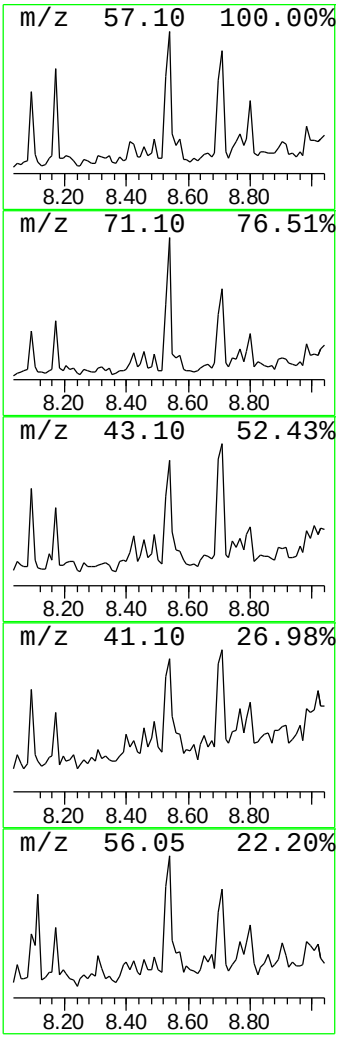
****

Peak Number 5 Octane, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.54	4.33 ng	131177	Naphthalene-d8	8.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	83
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	78
3	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	72
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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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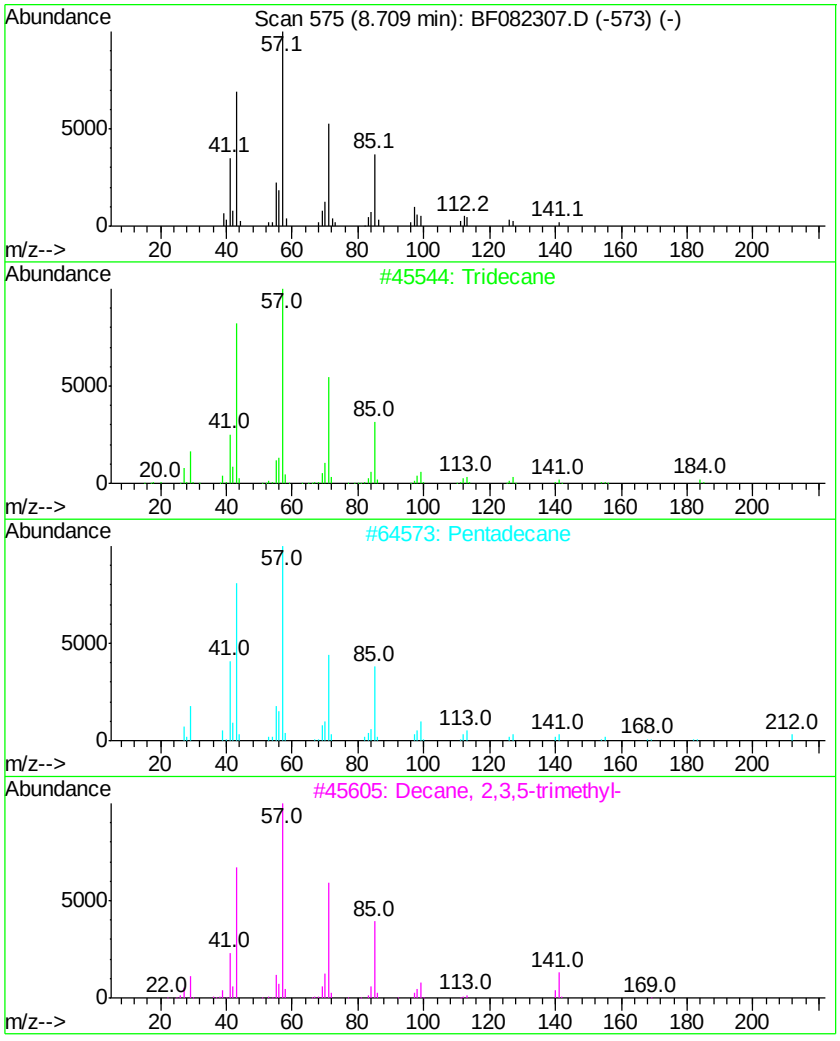
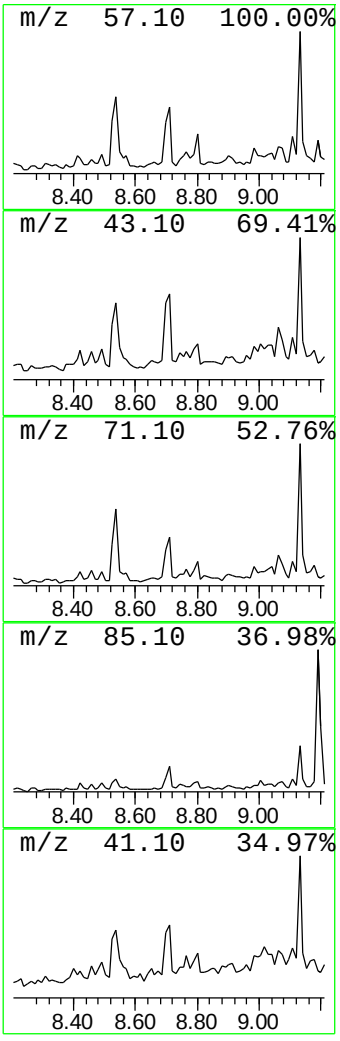
Peak Number 6 Tridecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	2.73 ng	82845	Naphthalene-d8	8.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	72
2	Pentadecane	212	C15H32	000629-62-9	64
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	64
4	Undecane	156	C11H24	001120-21-4	64
5	Hexadecane	226	C16H34	000544-76-3	64





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

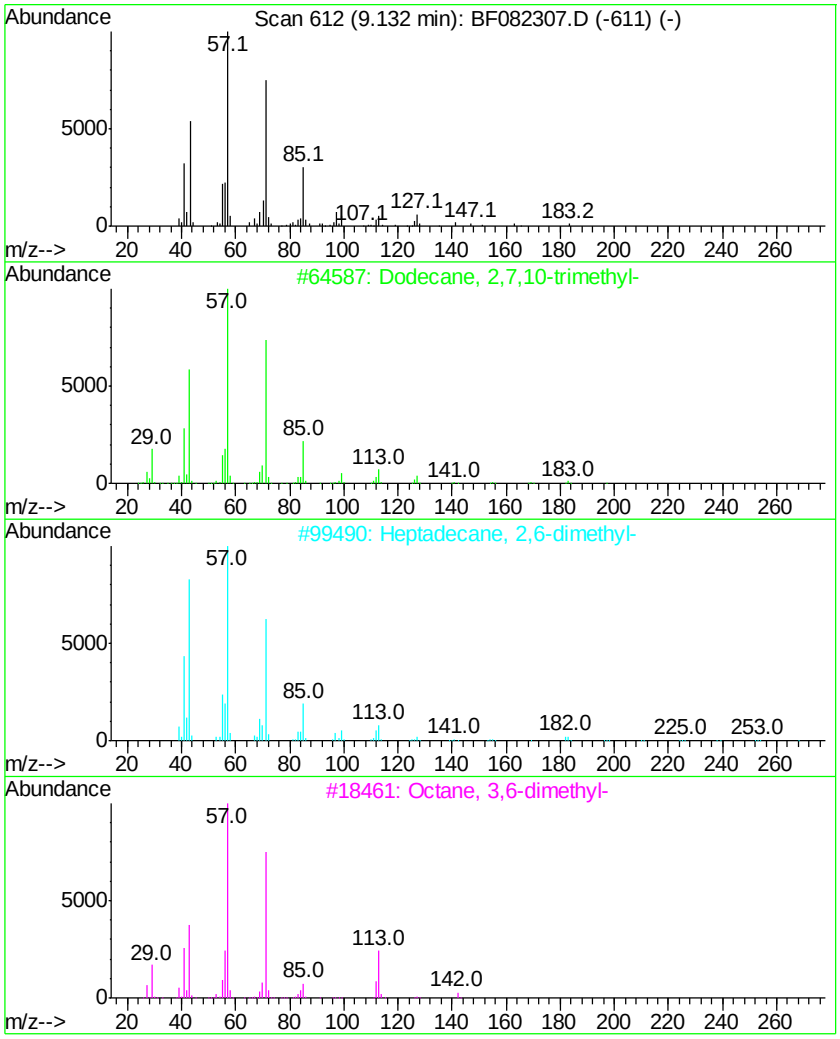
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

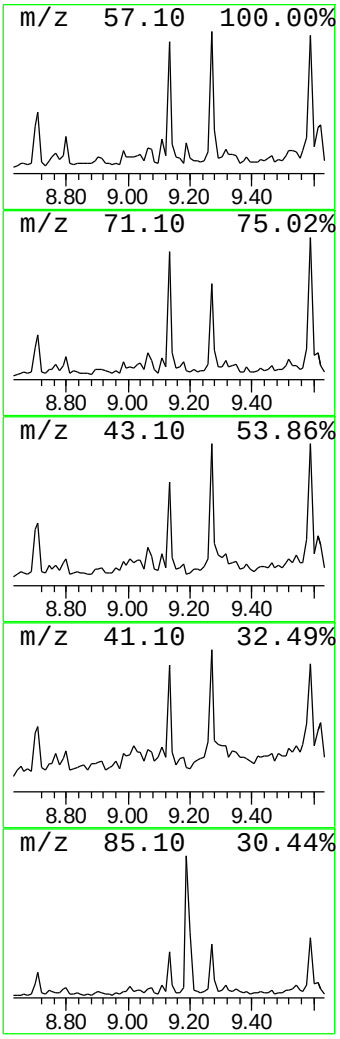
****

Peak Number 7 Dodecane, 2,7,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.13	3.27 ng	117350	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
2	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	72
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
4	Undecane, 6-ethyl-	184	C13H28	017312-60-6	64
5	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	64





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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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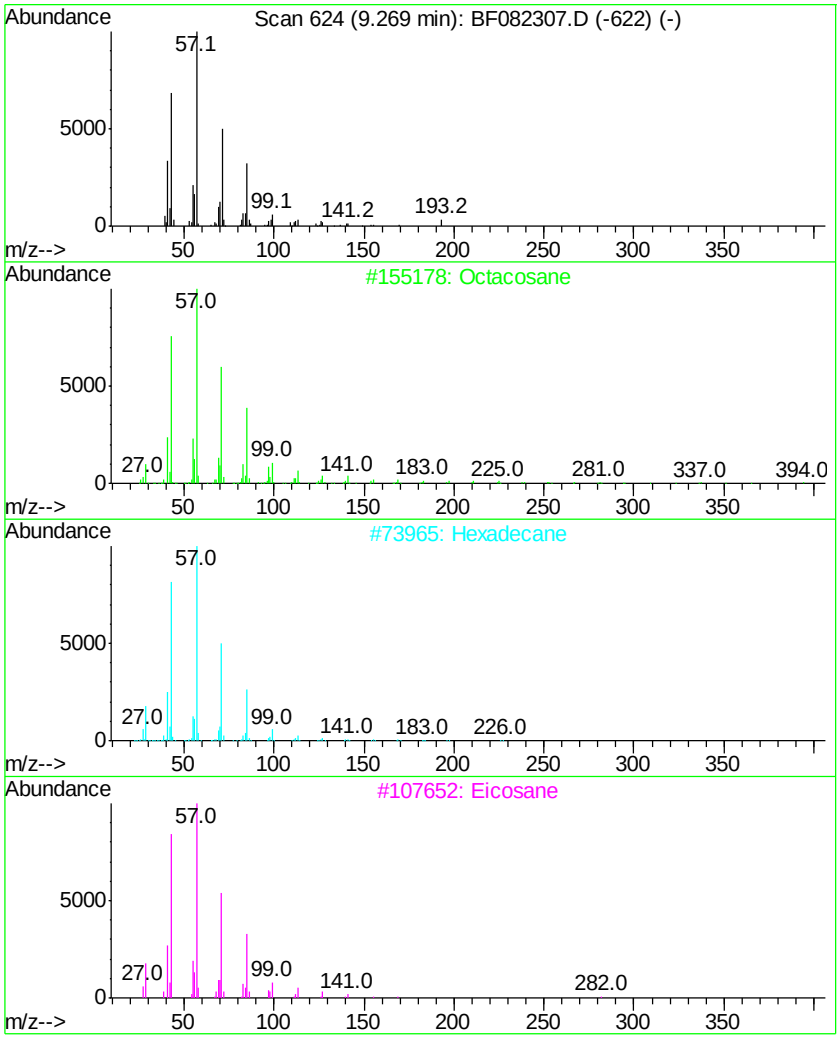
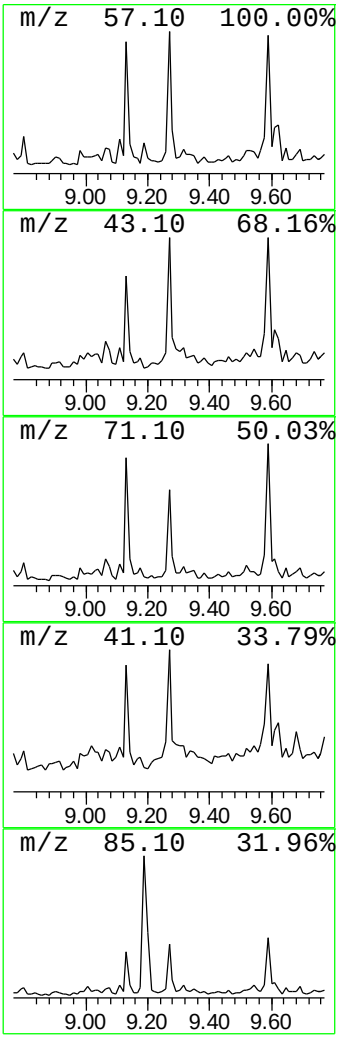
Peak Number 8 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.27	4.49 ng	161262	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	90
2	Hexadecane	226	C16H34	000544-76-3	90
3	Eicosane	282	C20H42	000112-95-8	87
4	Tetradecane	198	C14H30	000629-59-4	87
5	Pentadecane	212	C15H32	000629-62-9	86





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

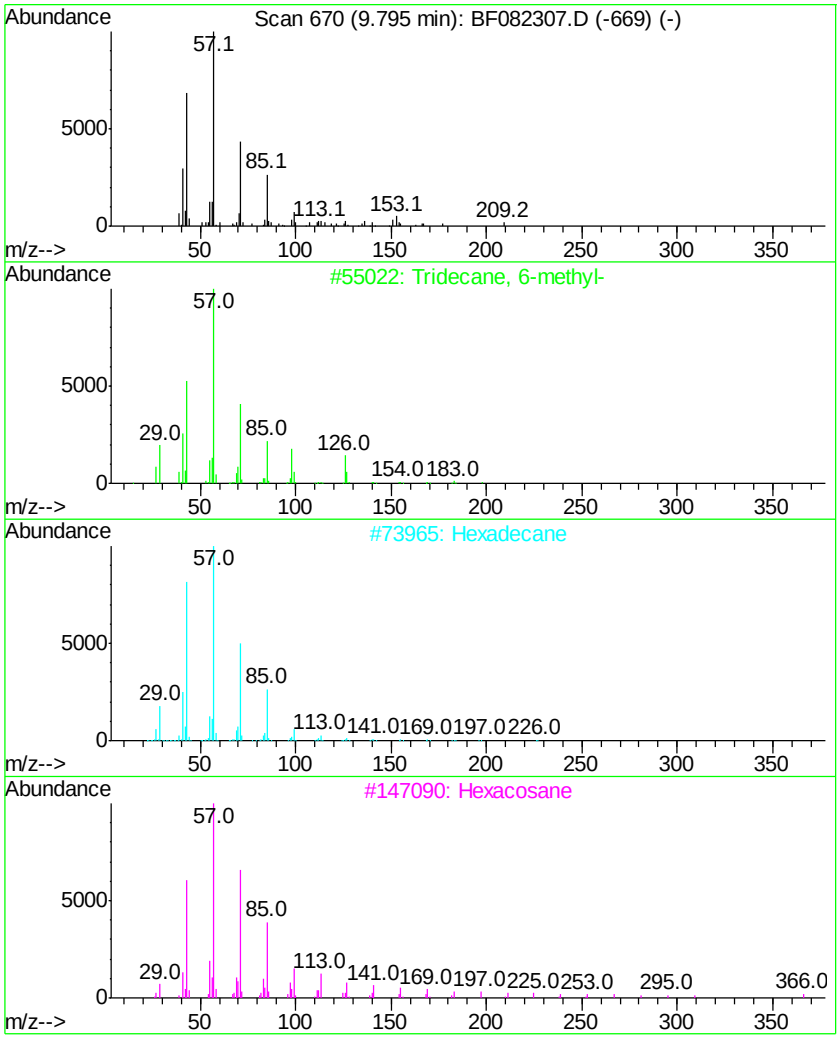
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

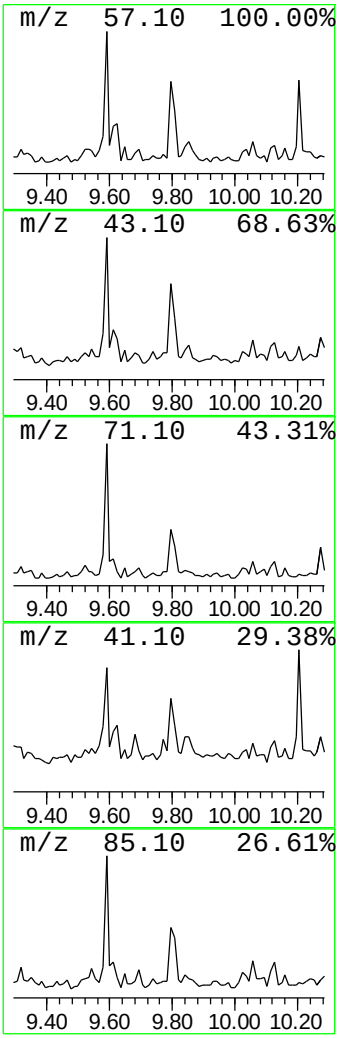
****

Peak Number 9 Tridecane, 6-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.79	3.11 ng	111735	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 6-methyl-	198	C14H30	013287-21-3	72
2	Hexadecane	226	C16H34	000544-76-3	72
3	Hexacosane	366	C26H54	000630-01-3	72
4	Octacosane	394	C28H58	000630-02-4	72
5	Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	59





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Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

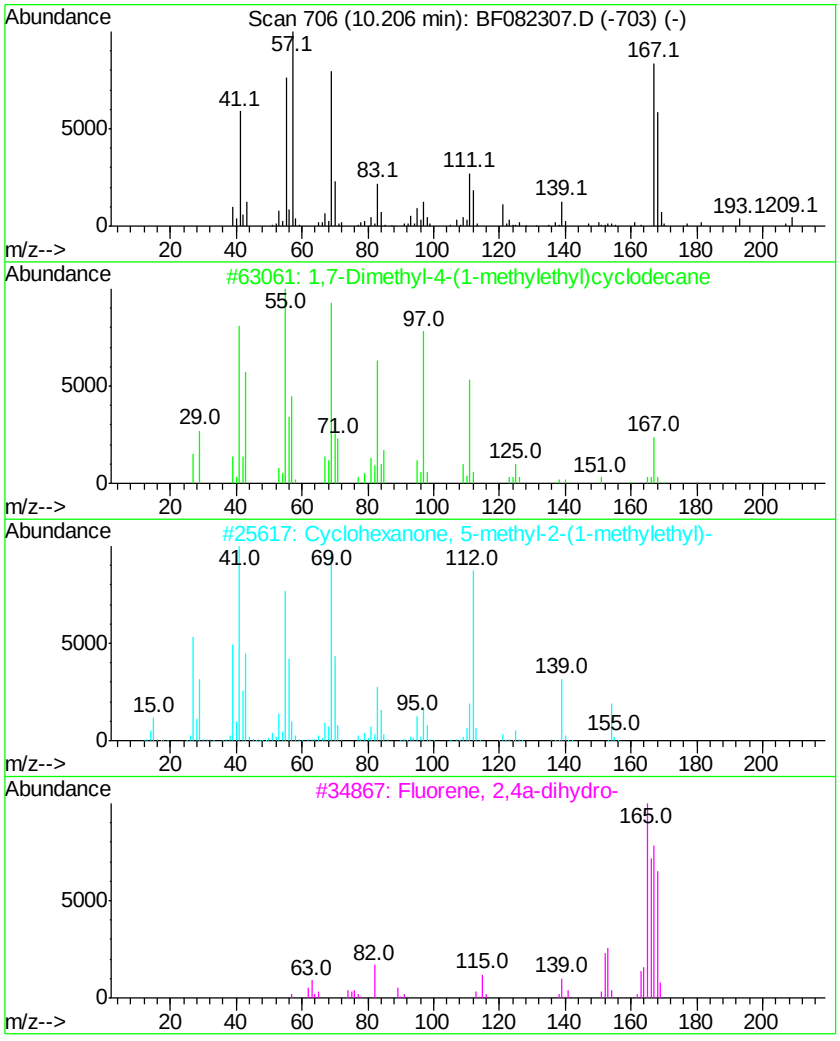
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

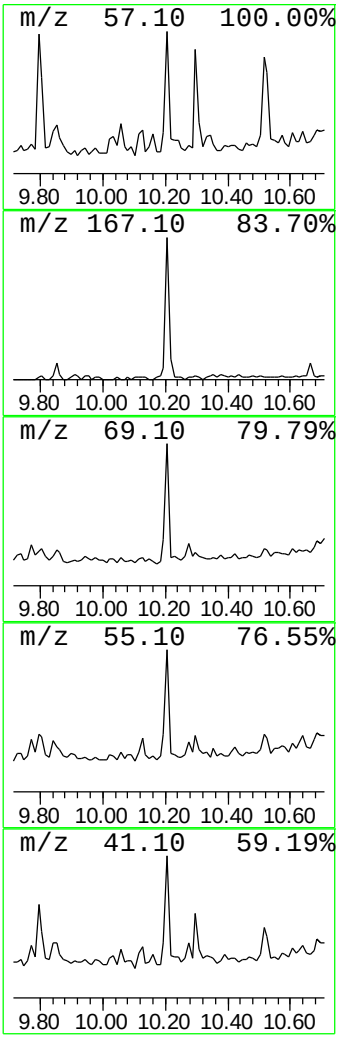
****

Peak Number 10 unknown10.21 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.21	4.93 ng	176986	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,7-Dimethyl-4-(1-methylethyl)cy...	210	C15H30	000645-10-3	38
2	Cyclohexanone, 5-methyl-2-(1-met...	154	C10H18O	010458-14-7	35
3	Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	30
4	Acetamide, N-cycloheptyl-2,2-dip...	307	C21H25NO	351062-01-6	27
5	Heptane, 4-methylene-	112	C8H16	015918-08-8	27





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

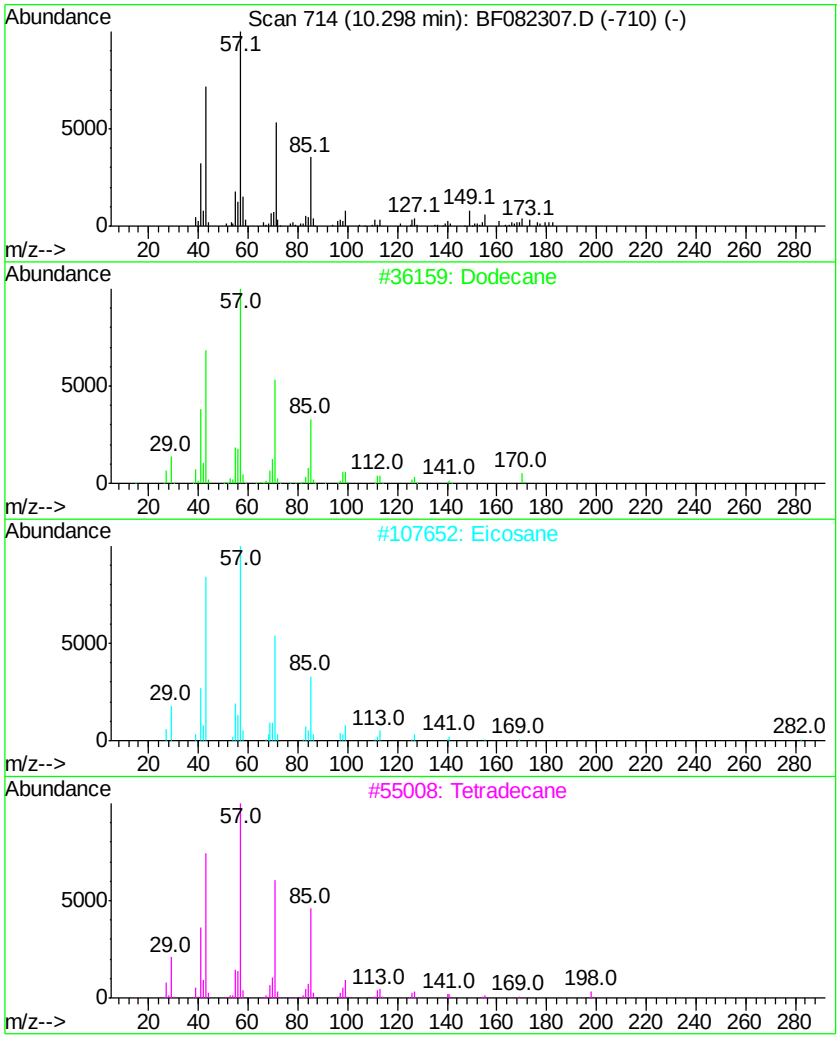
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

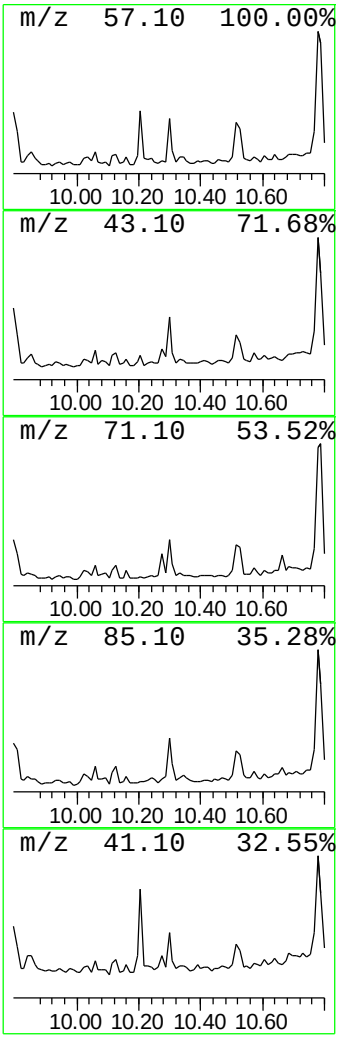
****

Peak Number 11 Dodecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	3.57 ng	128124	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	72
2	Eicosane	282	C20H42	000112-95-8	64
3	Tetradecane	198	C14H30	000629-59-4	64
4	Heptadecane	240	C17H36	000629-78-7	59
5	Hexadecane	226	C16H34	000544-76-3	59





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

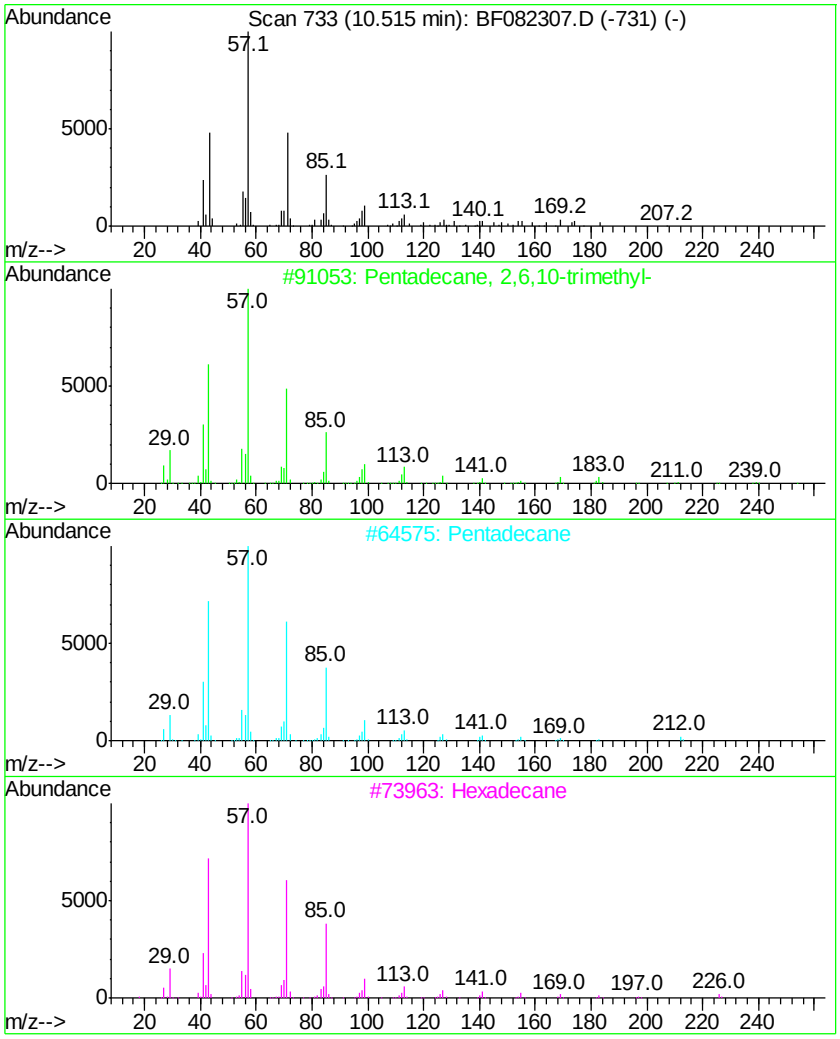
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

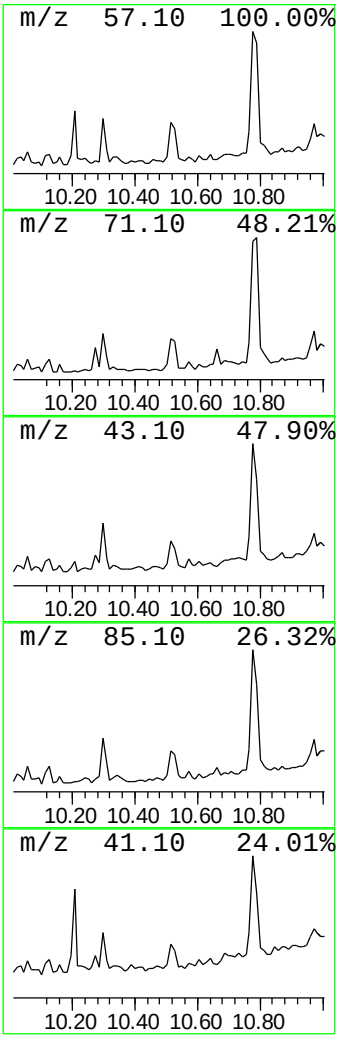
****

Peak Number 12 Pentadecane, 2,6,10-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	2.86 ng	102466	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	90
2	Pentadecane	212	C15H32	000629-62-9	86
3	Hexadecane	226	C16H34	000544-76-3	86
4	Eicosane	282	C20H42	000112-95-8	86
5	Heptadecane	240	C17H36	000629-78-7	80





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Acq On : 15 Oct 2015 14:13
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Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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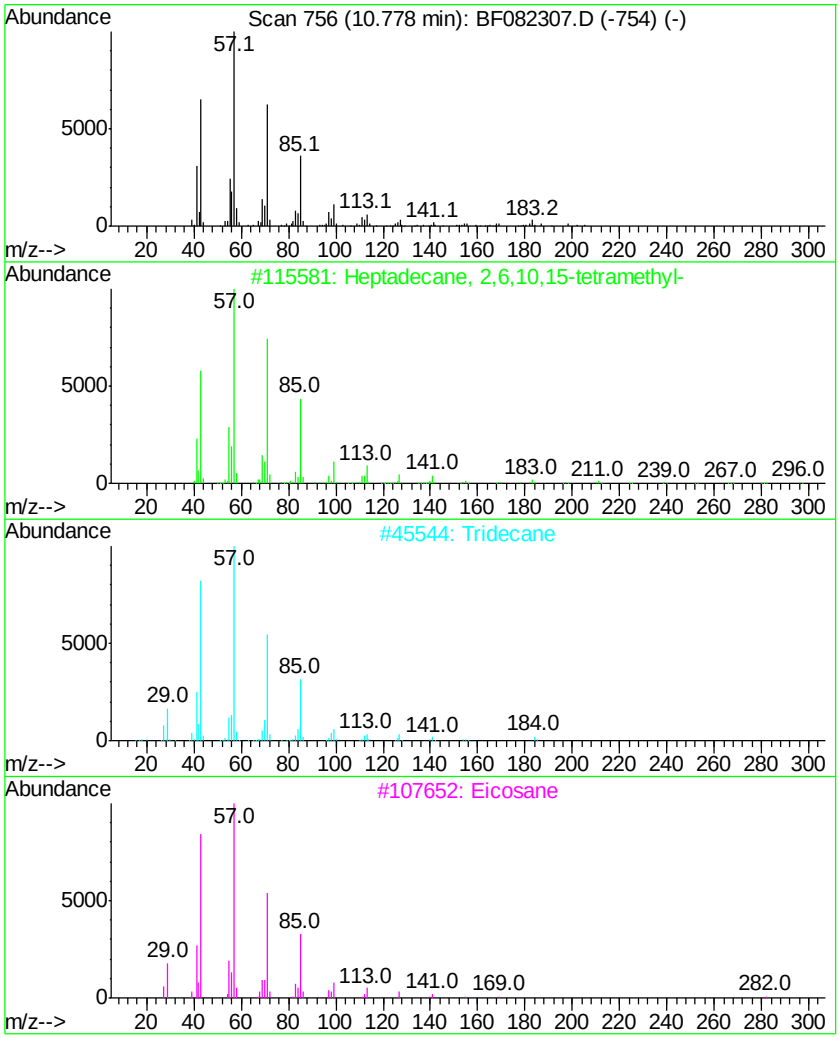
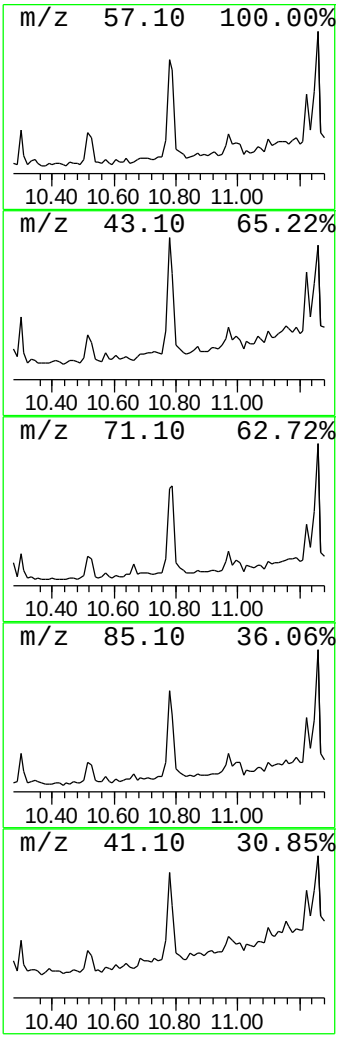
Peak Number 13 Heptadecane, 2,6,10,15-tetr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	11.72 ng	376719	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2	Tridecane	184	C13H28	000629-50-5	90
3	Eicosane	282	C20H42	000112-95-8	87
4	Heptadecane	240	C17H36	000629-78-7	86
5	Tetradecane	198	C14H30	000629-59-4	80





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Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
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Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
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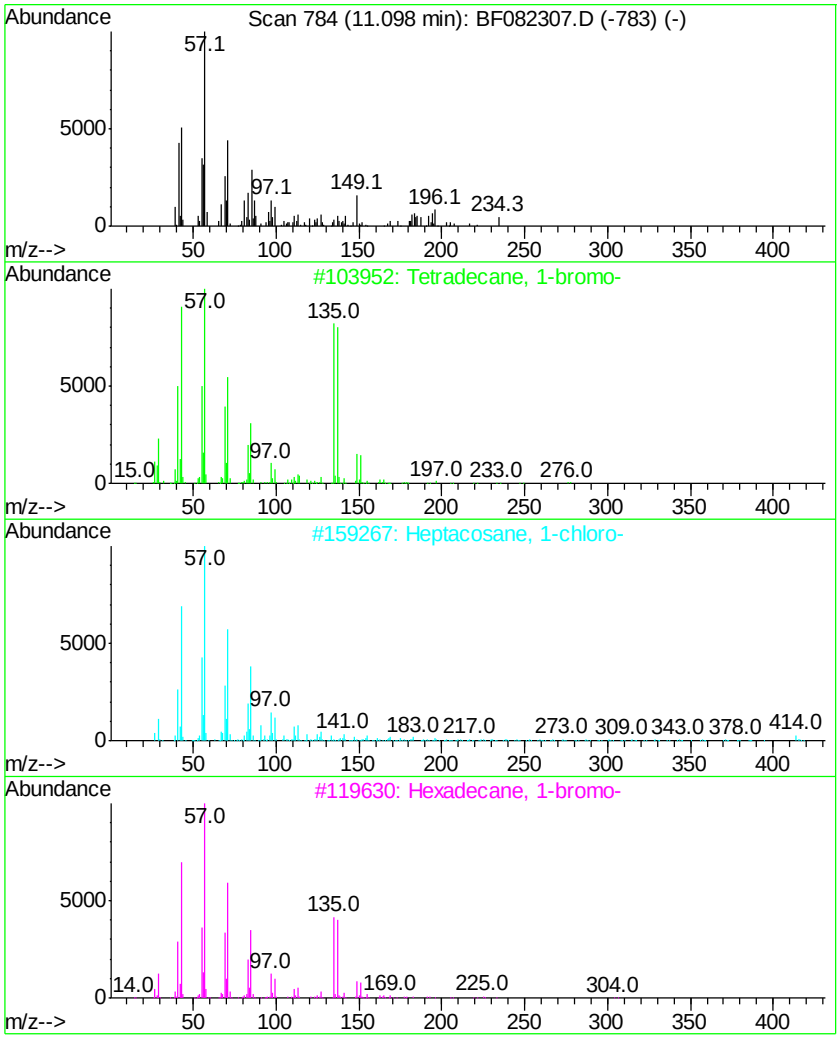
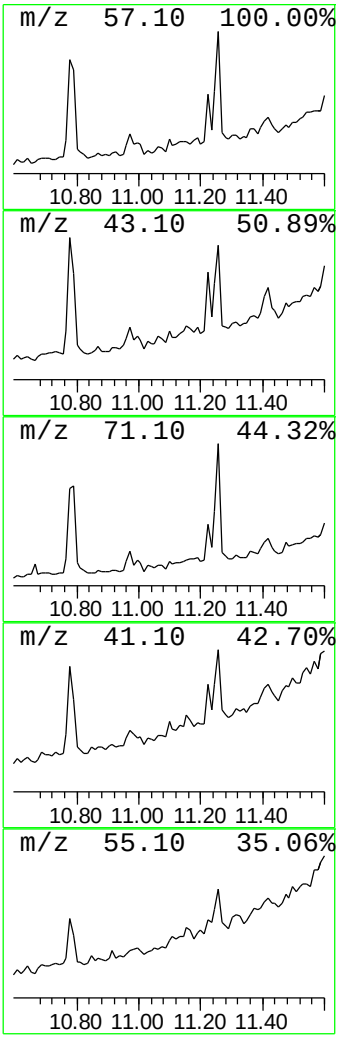
Peak Number 15 Tetradecane, 1-bromo- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.34 ng	75166	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	59
2	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	58
3	Hexadecane, 1-bromo-	304	C16H33Br	000112-82-3	53
4	Tetratetracontane	619	C44H90	007098-22-8	53
5	Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	53





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
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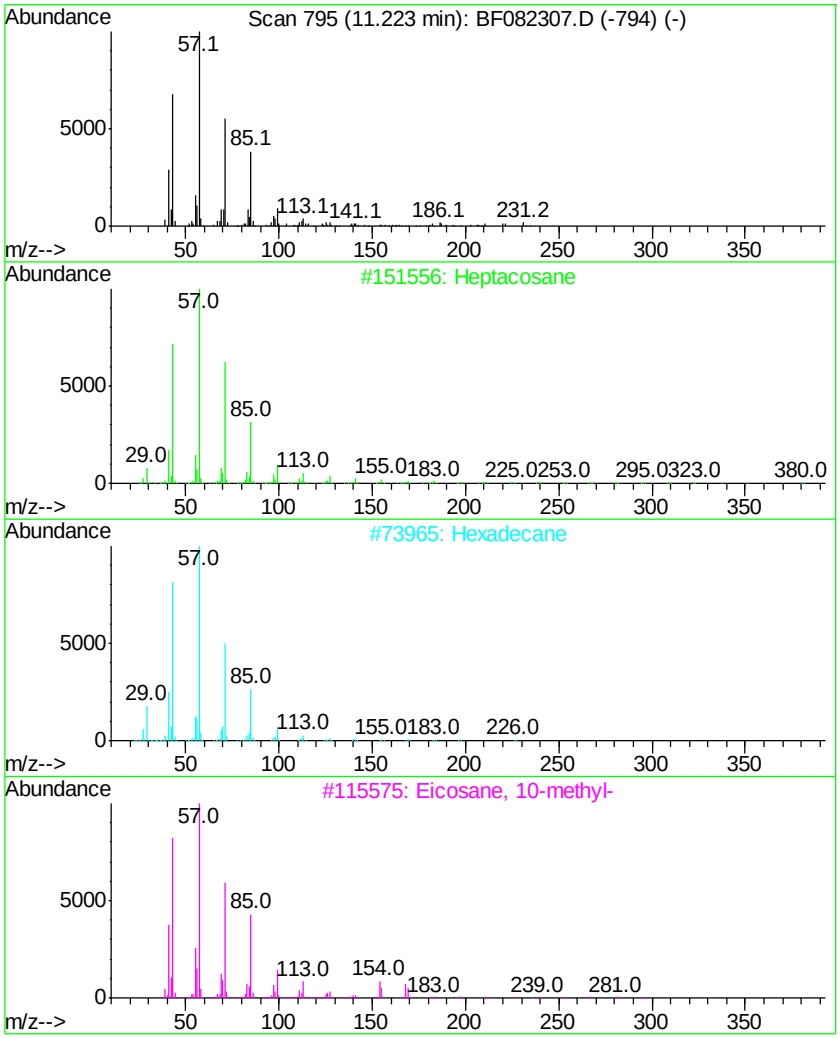
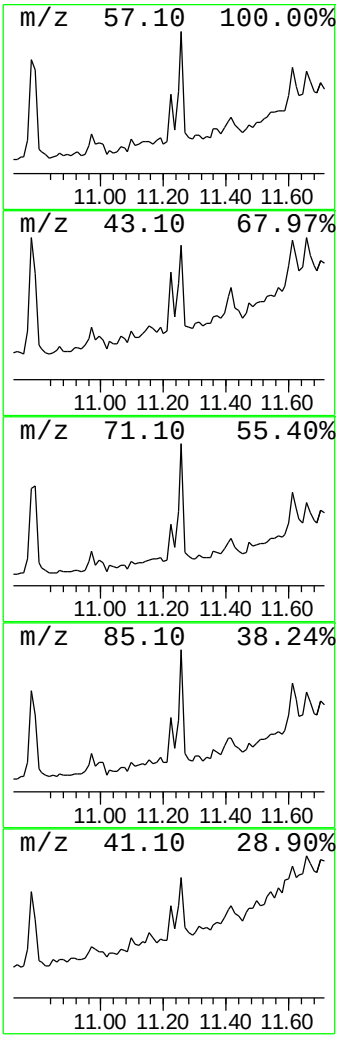
Peak Number 16 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	3.28 ng	105511	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	80
2	Hexadecane	226	C16H34	000544-76-3	64
3	Eicosane, 10-methyl-	296	C21H44	054833-23-7	64
4	Heneicosane	296	C21H44	000629-94-7	59
5	Tridecane	184	C13H28	000629-50-5	47





Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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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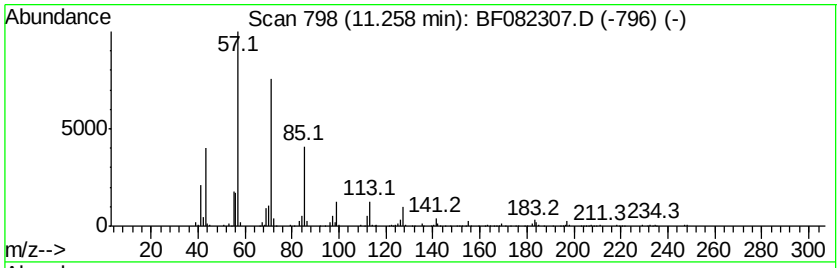
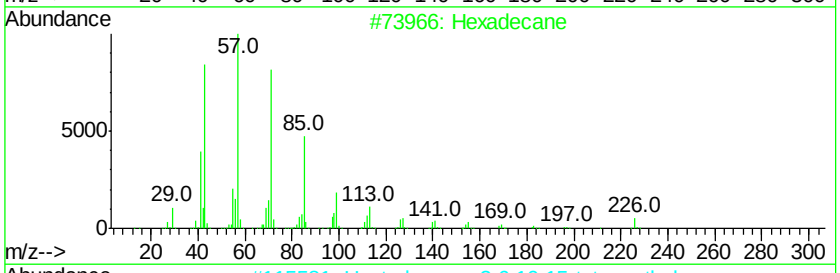
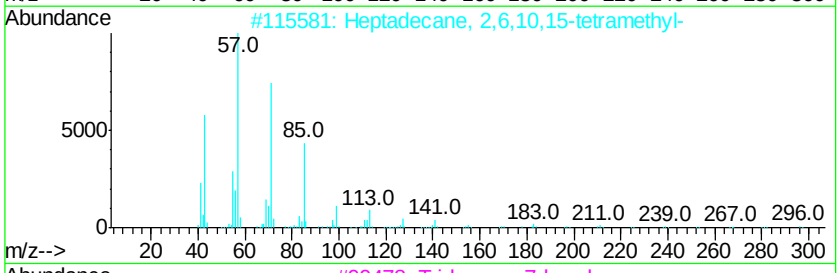
Peak Number 17 Hexadecane Concentration Rank 5

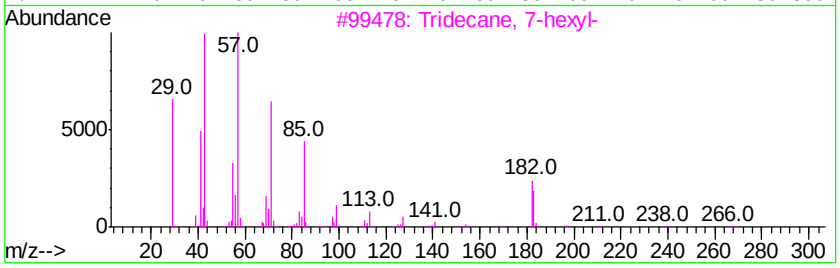
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	8.74 ng	281091	Phenanthrene-d10	11.36

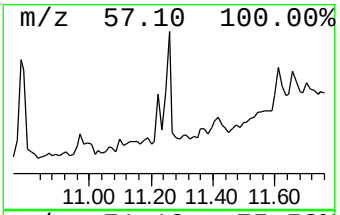
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	90
4	Heptadecane	240	C17H36	000629-78-7	90
5	Tetratetracontane	619	C44H90	007098-22-8	86

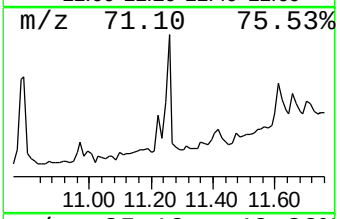


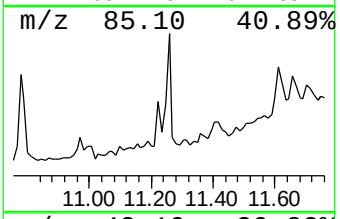


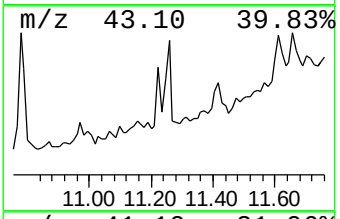


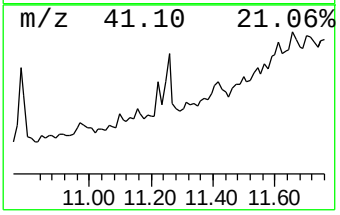












Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082307.D
Acq On : 15 Oct 2015 14:13
Operator : UM/IZ
Sample : G4016-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CS-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
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-Pentanone, 4-hy...	4.99	43.9	ng	934341	1	6.82	425660	20.0
unknown6.59	6.59	73.0	ng	1554370	1	6.82	425660	20.0
Benzene, 1,2,3-tr...	6.64	5.2	ng	110758	1	6.82	425660	20.0
Octane, 2,6-dimet...	8.54	4.3	ng	131177	2	8.11	606407	20.0
Tridecane	8.71	2.7	ng	82845	2	8.11	606407	20.0
Dodecane, 2,7,10-...	9.13	3.3	ng	117350	3	9.87	717663	20.0
Octacosane	9.27	4.5	ng	161262	3	9.87	717663	20.0
Tridecane, 6-methyl-	9.79	3.1	ng	111735	3	9.87	717663	20.0
unknown10.21	10.21	4.9	ng	176986	3	9.87	717663	20.0
Dodecane	10.30	3.6	ng	128124	3	9.87	717663	20.0
Pentadecane, 2,6,...	10.51	2.9	ng	102466	3	9.87	717663	20.0
Heptadecane, 2,6,...	10.78	11.7	ng	376719	4	11.36	642942	20.0
Tetradecane, 1-br...	11.10	2.3	ng	75166	4	11.36	642942	20.0
Heptacosane	11.22	3.3	ng	105511	4	11.36	642942	20.0
Hexadecane	11.26	8.7	ng	281091	4	11.36	642942	20.0