LSC Area Percent Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

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-BF112115.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	3.324	114	117	130	rVB	20169	73782	1.20%	0.159%
2	4.810	245	247	258	rBV	733532	957273	15.52%	2.062%
3	5.256	284	286	298	rBV	1554639	1750707	28.39%	3.771%
4	6.319	376	379	386	rBV	659549	1164586	18.88%	2.509%
5	6.422	386	388	391	rBV	3159092	3109149	50.42%	6.697%
6	6.650	406	408	411	rBV	1353068	1190586	19.31%	2.565%
7	6.753	415	417	419	rBV	132584	227701	3.69%	0.490%
8	6.810	419	422	424	rBV	4588414	4471192	72.50%	9.631%
9	7.222	455	458	461	rBV	3387491	3682007	59.70%	7.931%
10	7.462	476	479	481	rBV3	50412	85412	1.38%	0.184%
11	7.679	495	498	500	rBV2	81130	112643	1.83%	0.243%
12	7.782	504	507	509	rVB3	88482	116227	1.88%	0.250%
13	7.942	518	521	523	rBV	1579218	1626051	26.37%	3.503%
14	8.090	532	534	536	rBV2	46450	82011	1.33%	0.177%
15	8.228	544	546	548	rBV2	70352	80325	1.30%	0.173%
16	8.273	548	550	553	rBV	160137	163329	2.65%	0.352%
17	8.445	563	565	568	rVB2	111612	134208	2.18%	0.289%
18	8.536	568	573	576	rBV2	70237	149485	2.42%	0.322%
19	8.605	576	579	581	rVB2	74110	98010	1.59%	0.211%
20	8.765	589	593	595	rBV2	73652	123529	2.00%	0.266%
21	9.016	612	615	618	rVB	4850401	6167078	100.00%	13.284%
22	9.119	620	624	625	rBV3	56596	117748	1.91%	0.254%
23	9.153	625	627	629	rVB2	84847	116643	1.89%	0.251%
24	9.371	641	646	649	rBV	908301	1096577	17.78%	2.362%
25	9.416	649	650	652	rVV	184638	271591	4.40%	0.585%
26	9.462	652	654	657	rVV4	116021	222727	3.61%	0.480%
27	9.519	657	659	661	rVV2	87157	136541	2.21%	0.294%
28	9.599	665	666	668	rVV2	60461	104942	1.70%	0.226%
29	9.633	668	669	671	rVV	107122	109864	1.78%	0.237%
30	9.691	671	674	676	rVV	1541142	1795204	29.11%	3.867%
31	9.782	680	682	685	rVV3	39910	82371	1.34%	0.177%
32	9.828	685	686	688	rVV	88473	78493	1.27%	0.169%
33	9.896	688	692	693	rVV2	52461	101937	1.65%	0.220%
34	9.976	696	699	705	rVB2	69844	169907	2.76%	0.366%

LSC Area Percent Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
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Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

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-BF112115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.136	708	713	714	rBV2	93633	178349	2.89%	0.384%
36	10.171	714	716	720	rVB5	35527	92169	1.49%	0.199%
37	10.251	720	723	725	rBV4	66692	139000	2.25%	0.299%
38	10.308	725	728	730	rBV4	54753	124347	2.02%	0.268%
39	10.353	730	732	735	rBV	113232	143767	2.33%	0.310%
40	10.468	740	742	745	rBV2	2061217	2514968	40.78%	5.417%
41	10.559	747	750	752	rBV2	39978	64040	1.04%	0.138%
42	10.605	752	754	757	rBV	191619	284695	4.62%	0.613%
43	11.051	791	793	795	rBV	388377	391955	6.36%	0.844%
44	11.085	795	796	797	rVV	138334	101417	1.64%	0.218%
45	11.165	800	803	805	rBV	1532543	1586788	25.73%	3.418%
46	11.474	828	830	833	rBV	78861	113837	1.85%	0.245%
47	11.542	835	836	837	rVB	151075	107965	1.75%	0.233%
48	11.656	842	846	848	rBV5	77197	216649	3.51%	0.467%
49	11.714	849	851	855	rBV2	481801	690985	11.20%	1.488%
50	12.171	888	891	893	rBV	181037	330896	5.37%	0.713%
51	12.377	907	909	911	rBV2	175242	321618	5.22%	0.693%
52	12.502	917	920	923	rBV2	471096	852850	13.83%	1.837%
53	12.651	931	933	934	rBV	178611	280725	4.55%	0.605%
54	12.754	939	942	944	rVV	5045219	5280330	85.62%	11.374%
55	13.119	972	974	976	rBV	255655	422680	6.85%	0.910%
56	13.794	1030	1033	1035	rBV	928470	1392802	22.58%	3.000%
57	15.154	1150	1152	1155	rVB	696045	821088	13.31%	1.769%

Sum of corrected areas: 46423756

LSC Report - Integrated Chromatogram

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Data File : BF083136.D
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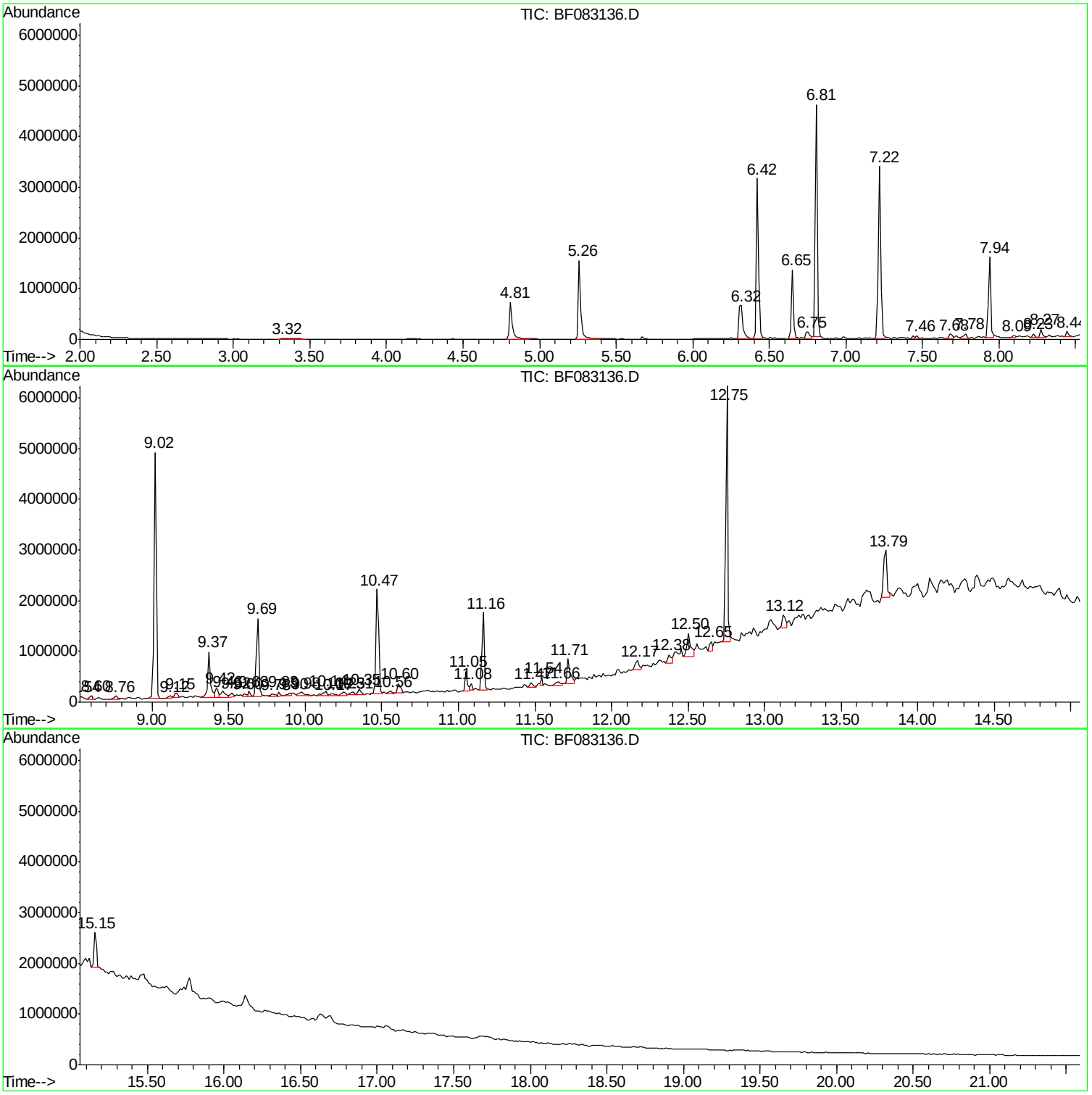
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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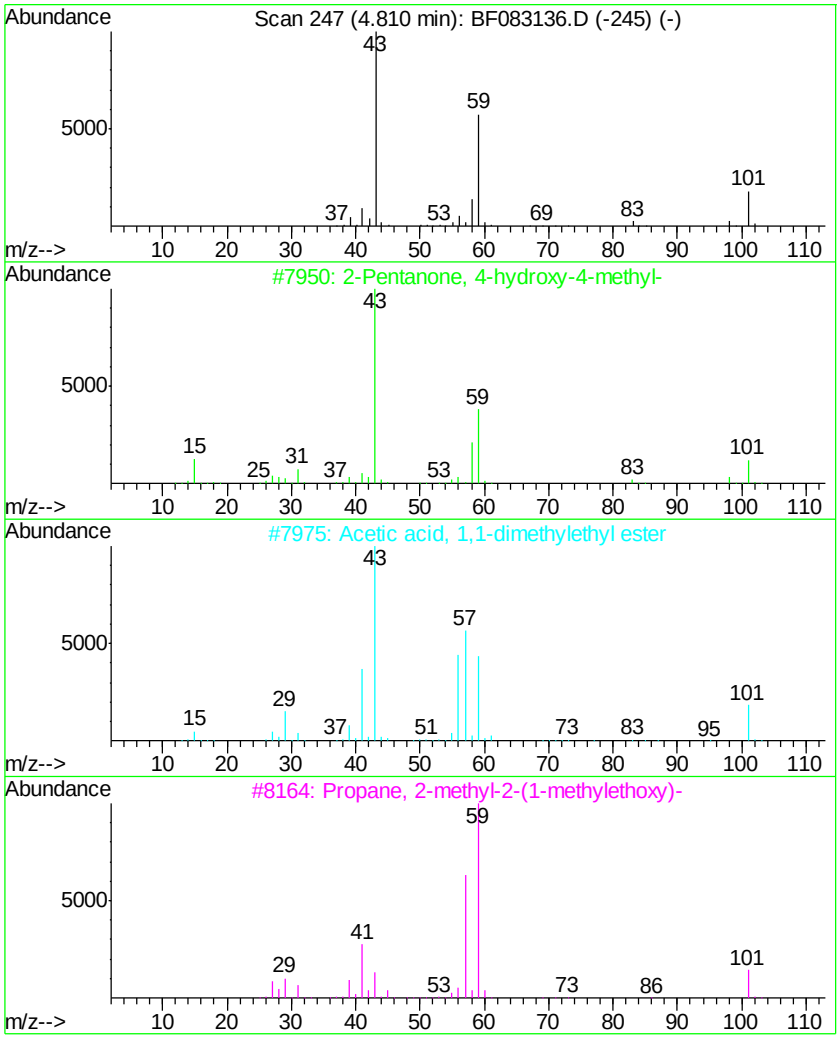
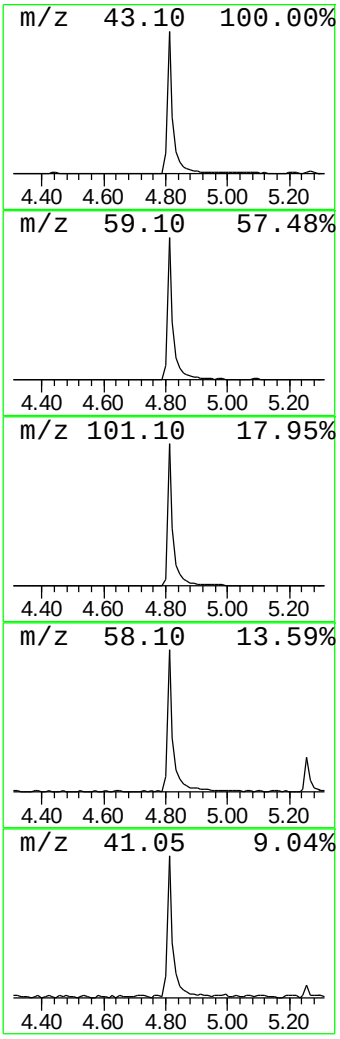
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.81	16.08 ng	957273	1,4-Dichlorobenzene-d4	6.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
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	36
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





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Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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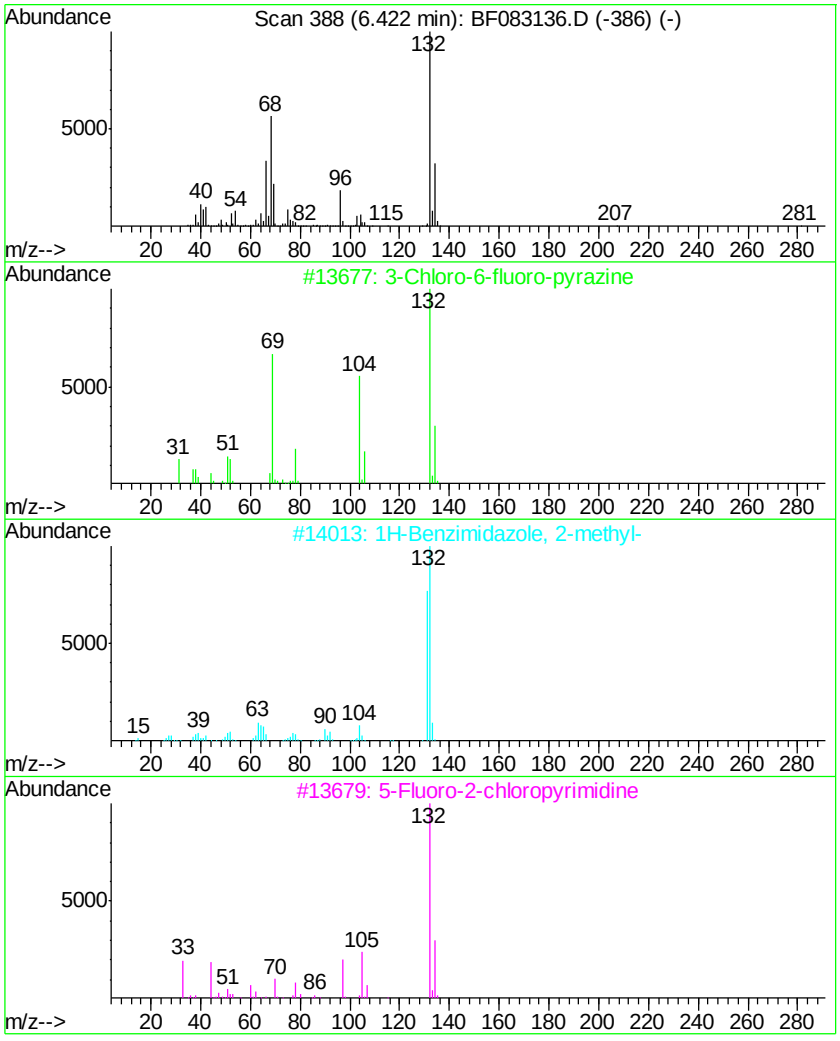
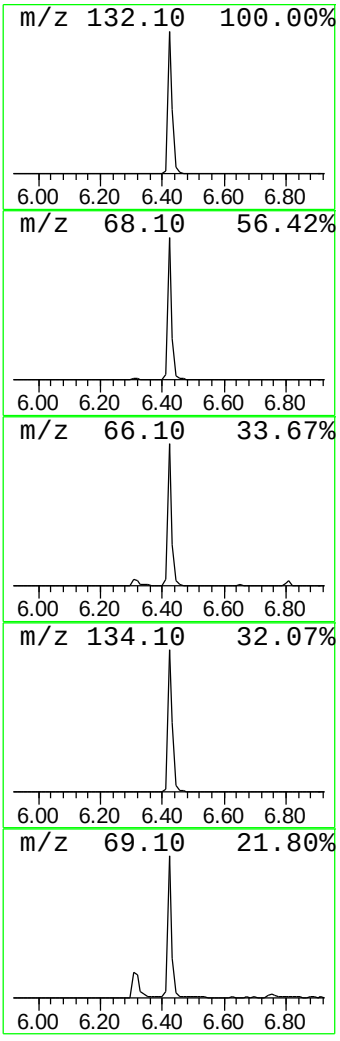
Peak Number 2 unknown6.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	52.23 ng	3109150	1,4-Dichlorobenzene-d4	6.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10





Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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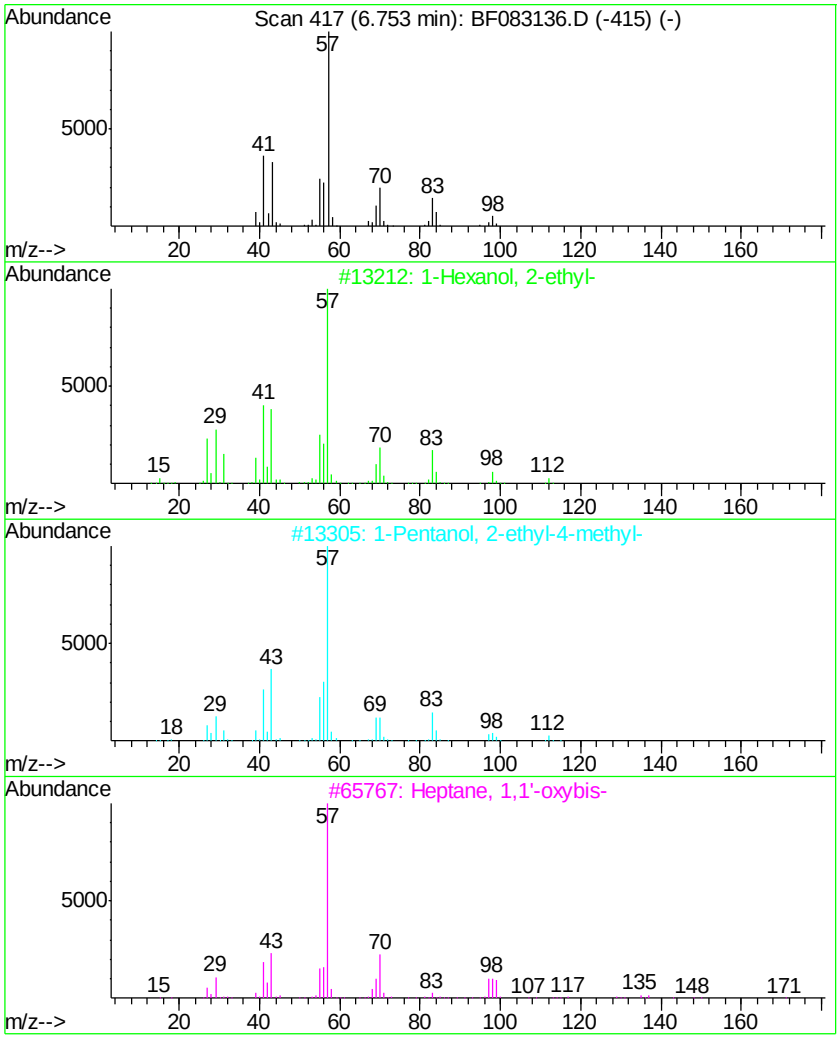
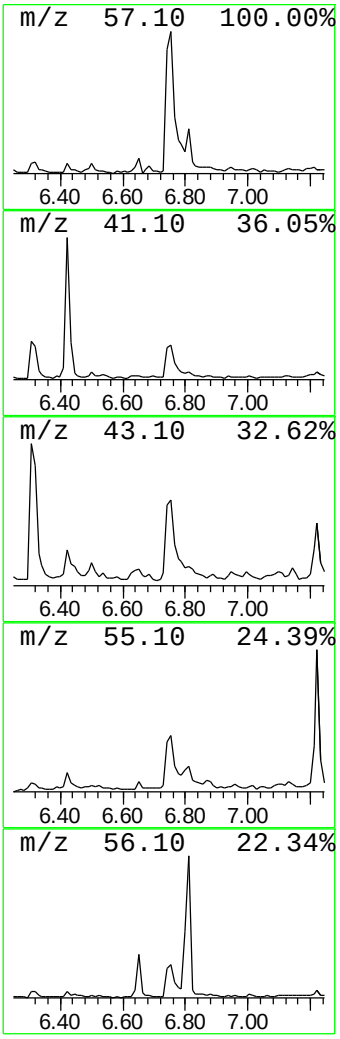
Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	3.83 ng	227701	1,4-Dichlorobenzene-d4	6.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2	1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	78
3	Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	53
4	Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	50
5	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	47





Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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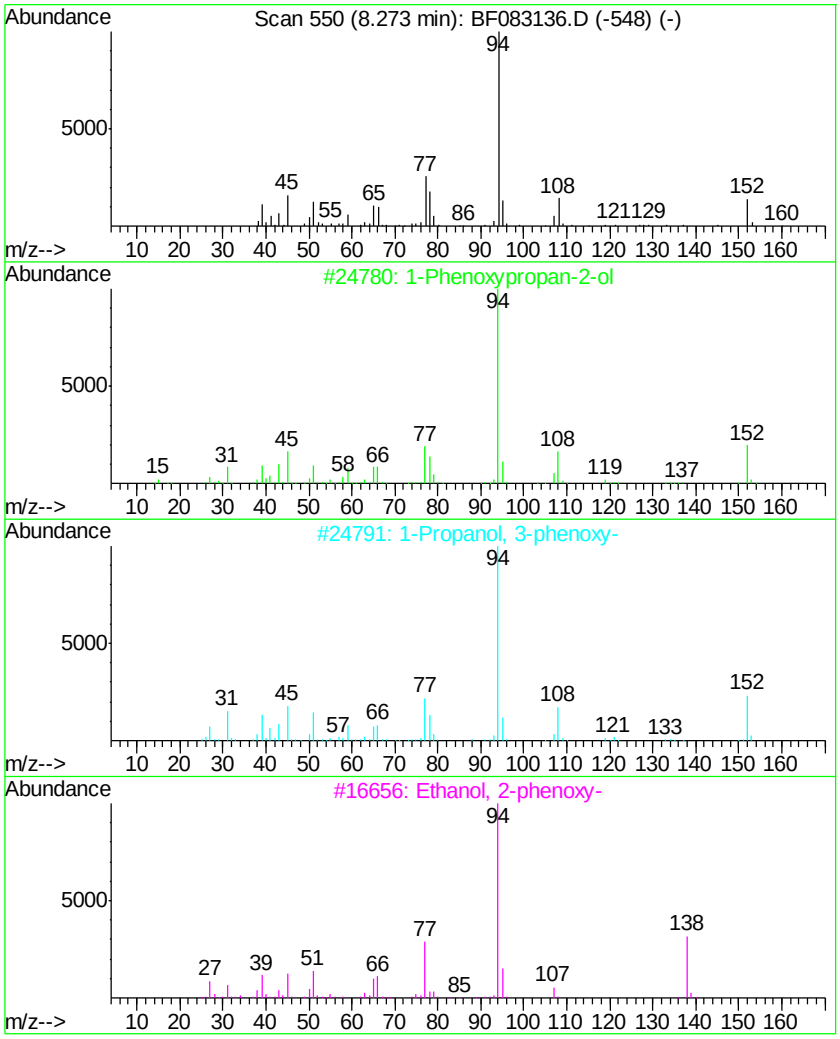
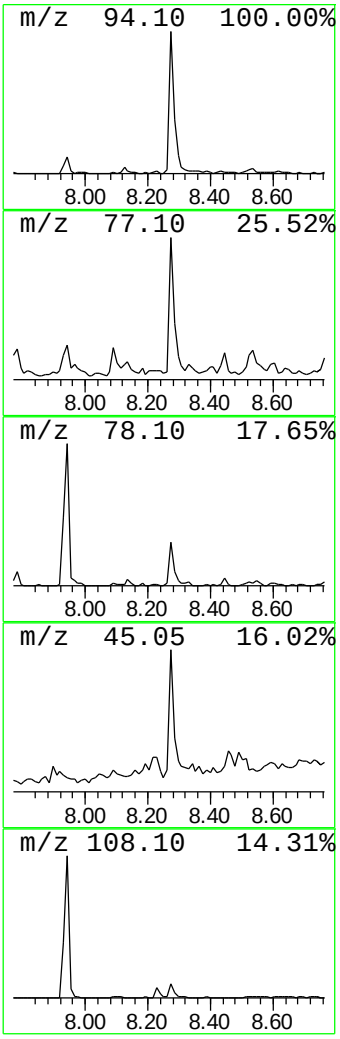
Peak Number 5 1-Phenoxypropan-2-ol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.27	2.01 ng	163329	Naphthalene-d8	7.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
2	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
3	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4	1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	59
5	Benzene, (1-methylpropoxy)-	150	C10H14O	010574-17-1	53





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Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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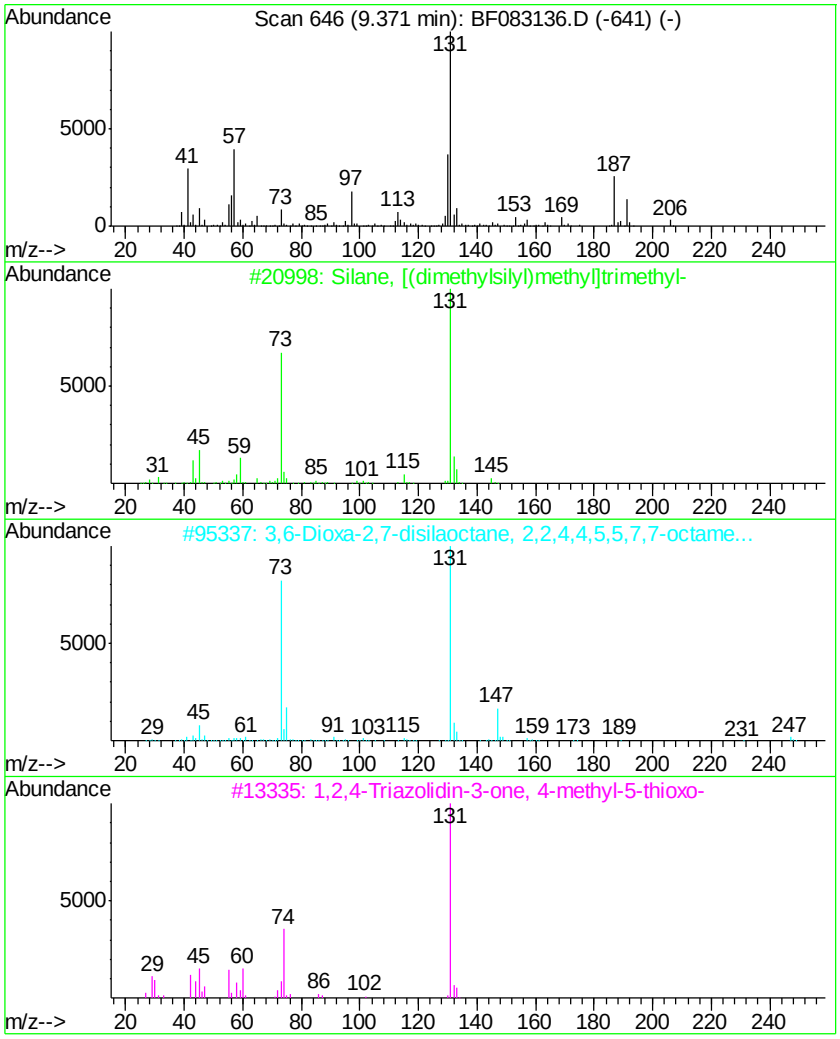
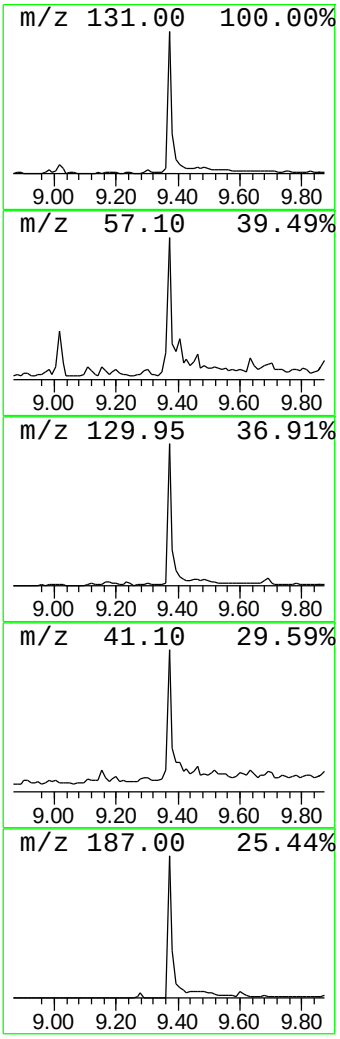
Peak Number 6 unknown9.37 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	12.22 ng	1096580	Acenaphthene-d10	9.69

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, [(dimethylsilyl)methyl...	146	C6H18Si2	001189-75-9	37
2	3,6-Dioxa-2,7-disilaoctane, 2,2,...	262	C12H30O2Si2	006730-96-7	25
3	1,2,4-Triazolidin-3-one, 4-methy...	131	C3H5N3OS	022244-61-7	25
4	1H-Indole, 4-methyl-	131	C9H9N	016096-32-5	23
5	1H-Indole, 5-methyl-	131	C9H9N	000614-96-0	9





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Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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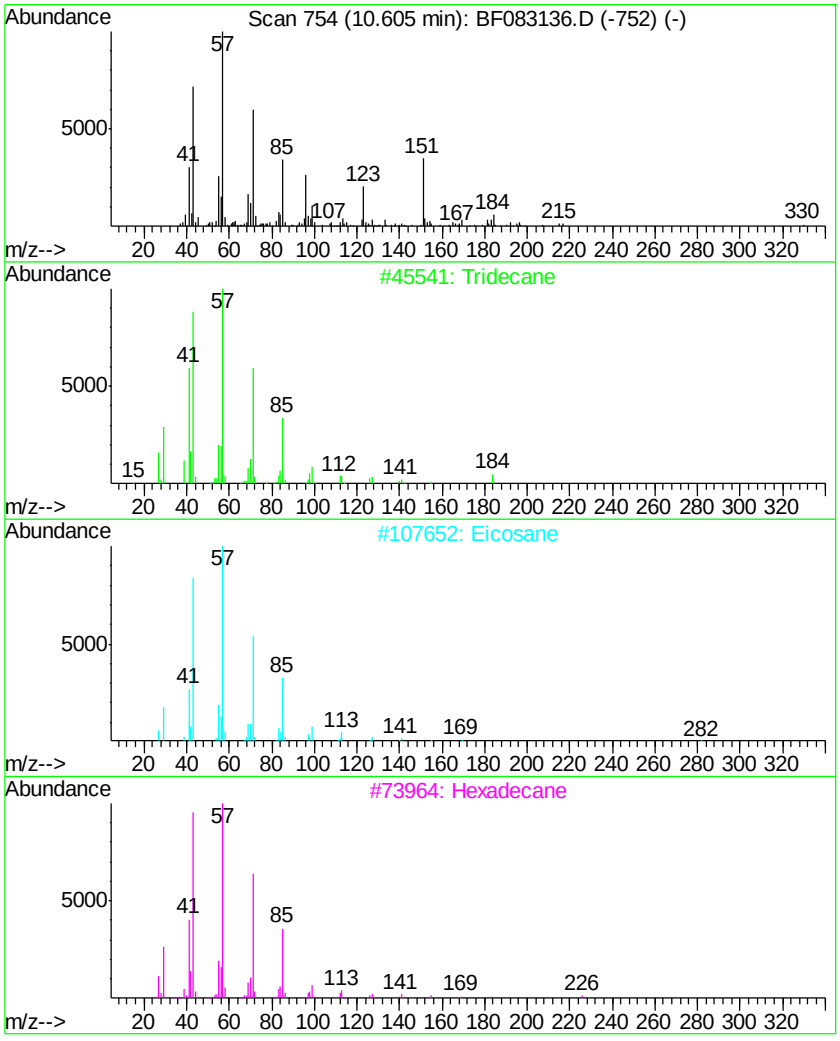
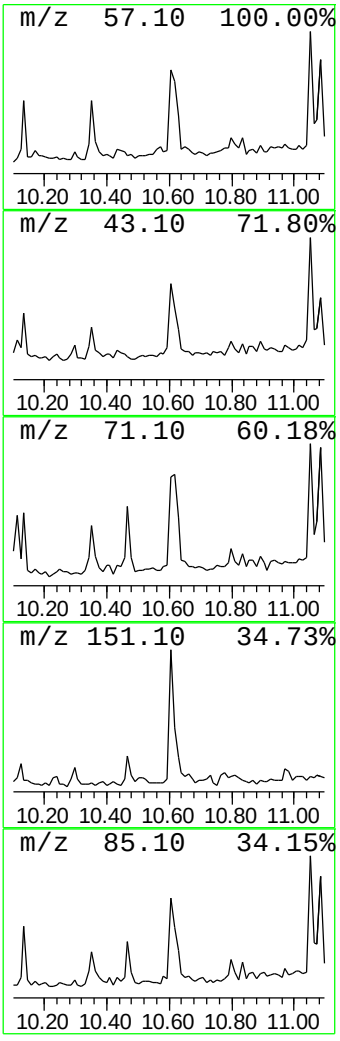
Peak Number 7 Tridecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.60	3.59 ng	284695	Phenanthrene-d10	11.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	78
2	Eicosane	282	C20H42	000112-95-8	49
3	Hexadecane	226	C16H34	000544-76-3	46
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	46
5	Tetradecane	198	C14H30	000629-59-4	46





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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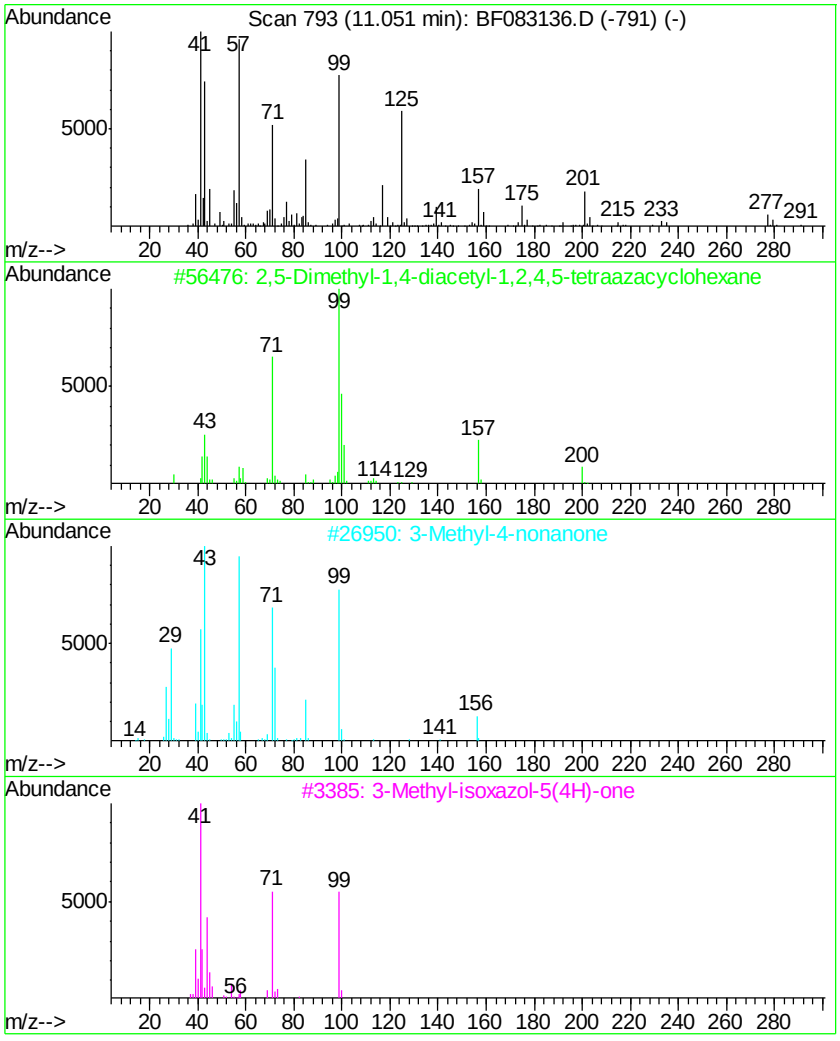
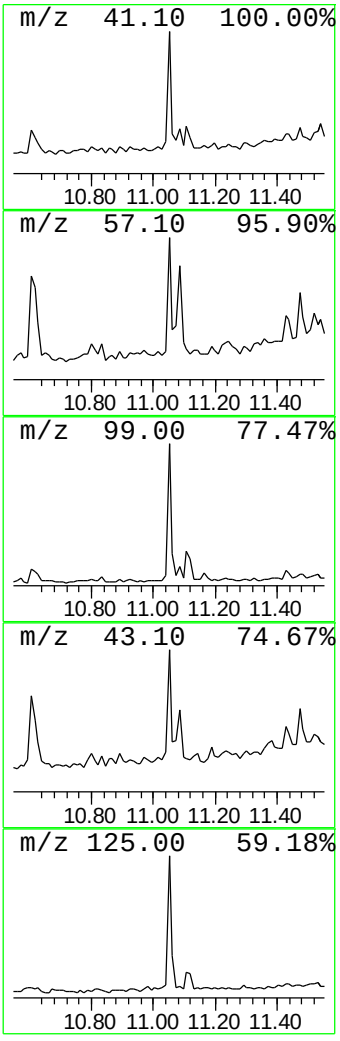
Peak Number 8 unknown11.05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.05	4.94 ng	391955	Phenanthrene-d10	11.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,5-Dimethyl-1,4-diacetyl-1,2,4,...	200	C8H16N4O2	021599-69-9	43
2	3-Methyl-4-nonanone	156	C10H20O	035778-39-3	38
3	3-Methyl-isoxazol-5(4H)-one	99	C4H5NO2	001517-96-0	35
4	3-Methyl-5-hydroxy-isoxazole	99	C4H5NO2	045469-93-0	35
5	Eicosane	282	C20H42	000112-95-8	15





Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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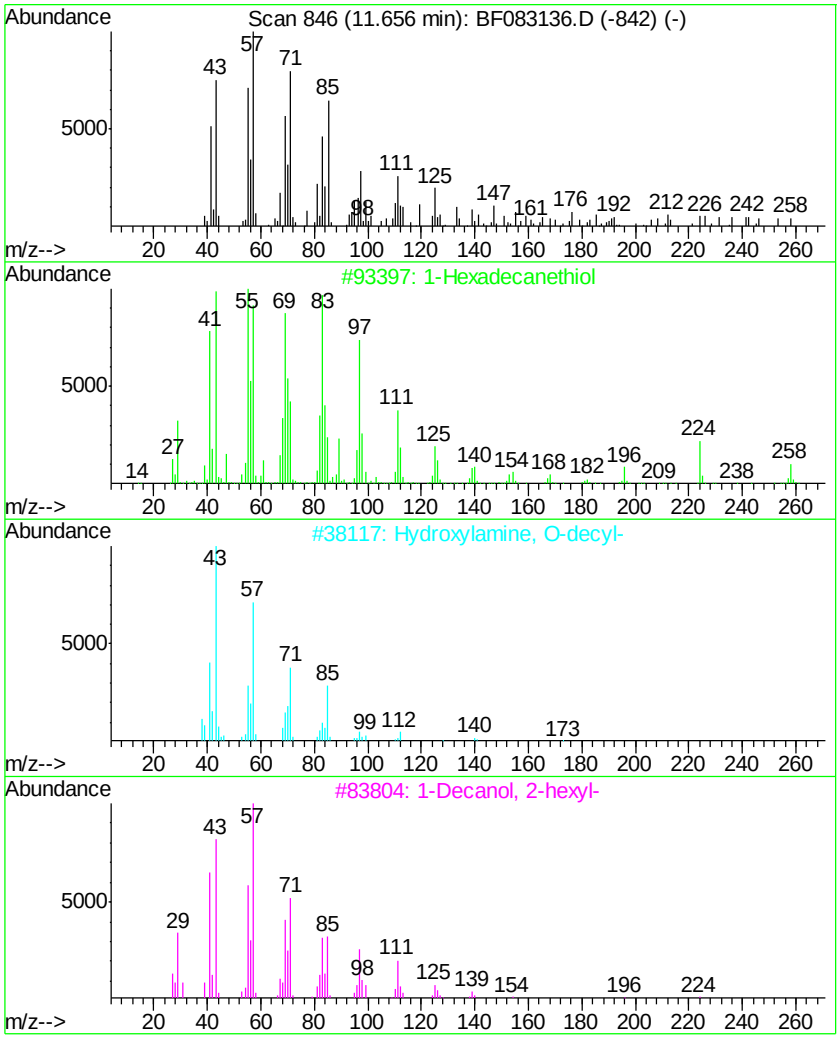
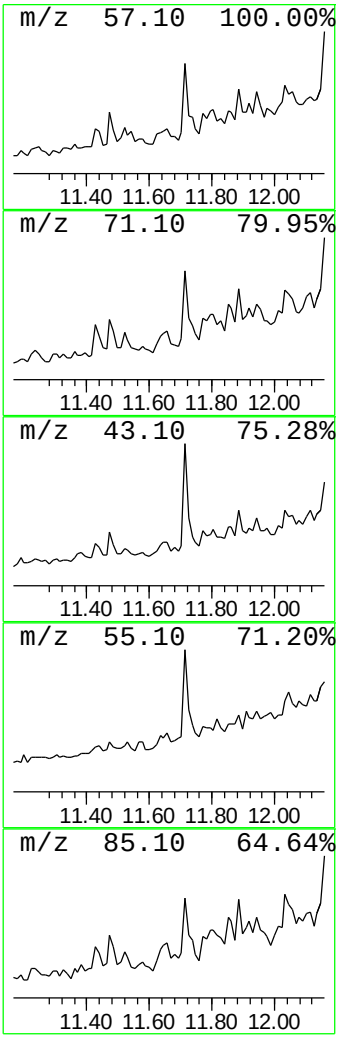
Peak Number 9 1-Hexadecanethiol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	2.73 ng	216649	Phenanthrene-d10	11.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanethiol	258	C16H34S	002917-26-2	64
2	Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	58
3	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	52
4	Hexadecane	226	C16H34	000544-76-3	50
5	1-Hexadecene	224	C16H32	000629-73-2	46





Library Search Compound Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

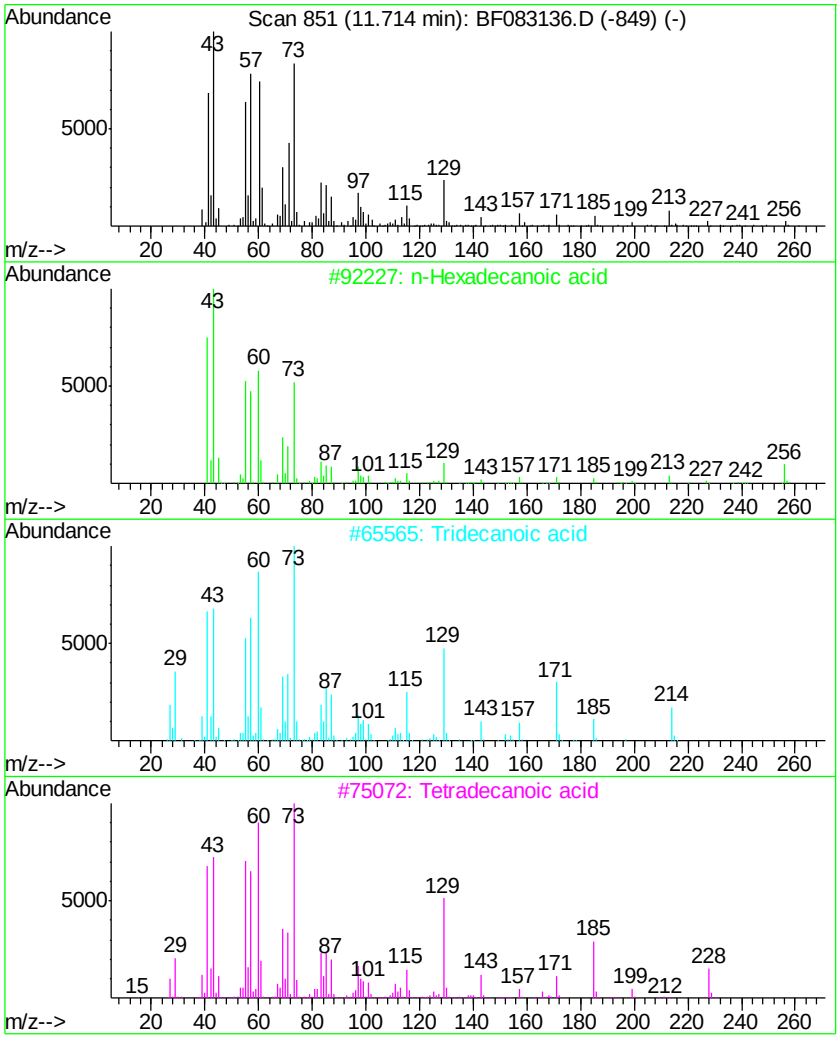
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

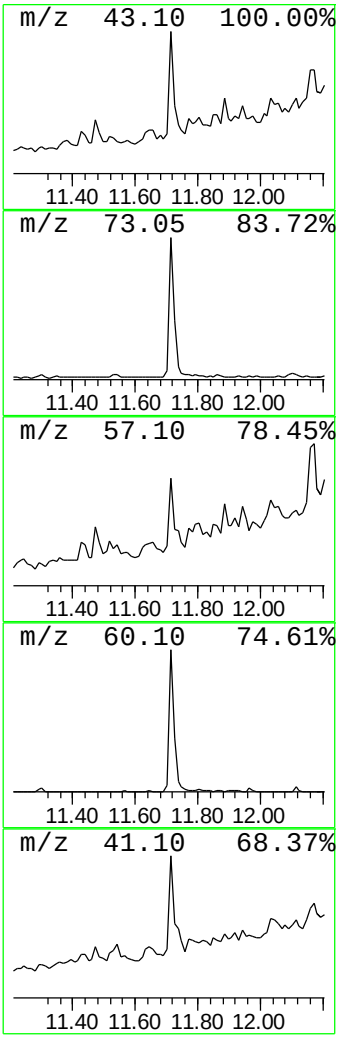
****

Peak Number 10 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	8.71 ng	690985	Phenanthrene-d10	11.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	Tridecanoic acid	214	C13H26O2	000638-53-9	89
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
4	Undecanoic acid	186	C11H22O2	000112-37-8	59
5	Amyl Nitrite	117	C5H11NO2	000110-46-3	38





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Sample : G4468-02
Misc :
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Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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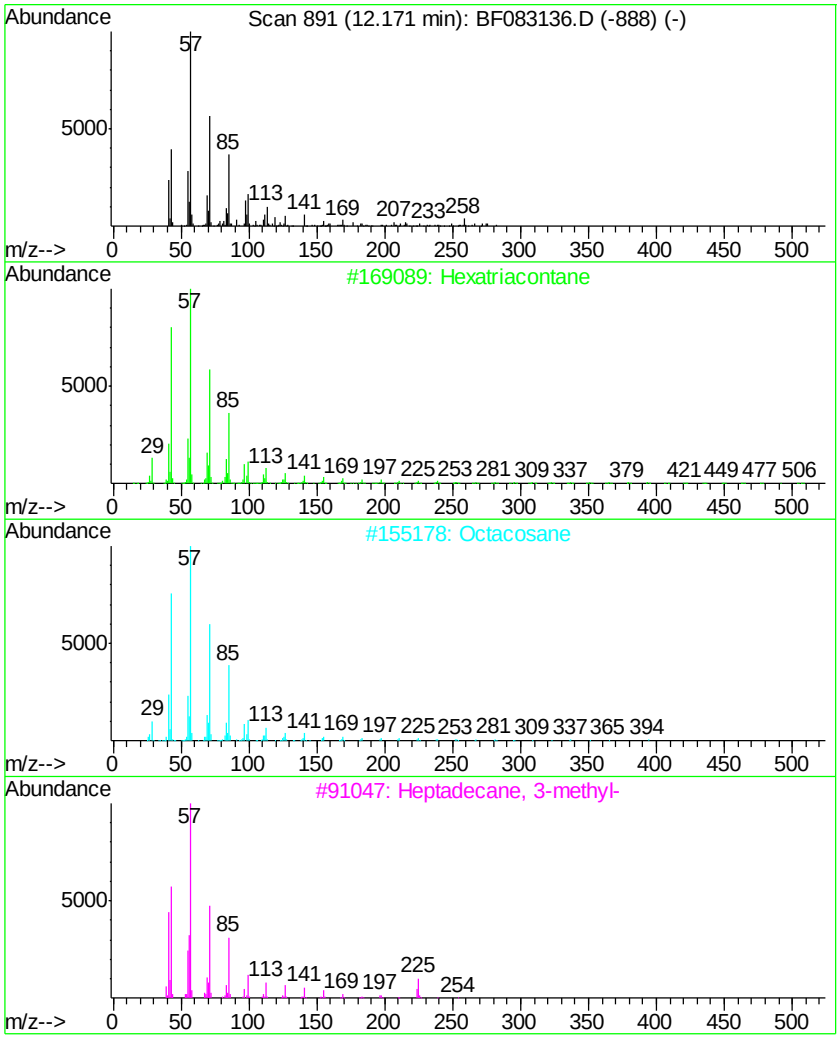
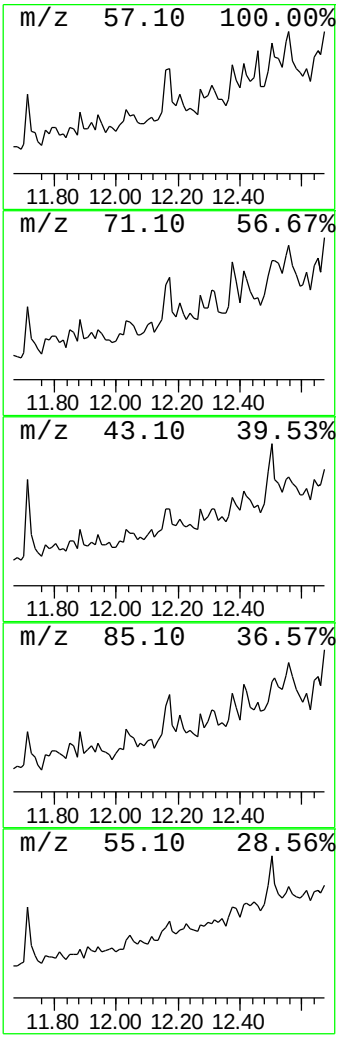
Peak Number 11 Hexatriacontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.17	4.17 ng	330896	Phenanthrene-d10	11.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexatriacontane	507	C36H74	000630-06-8	87
2	Octacosane	394	C28H58	000630-02-4	87
3	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	83
4	Octadecane	254	C18H38	000593-45-3	80
5	Heptadecane	240	C17H36	000629-78-7	80





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Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

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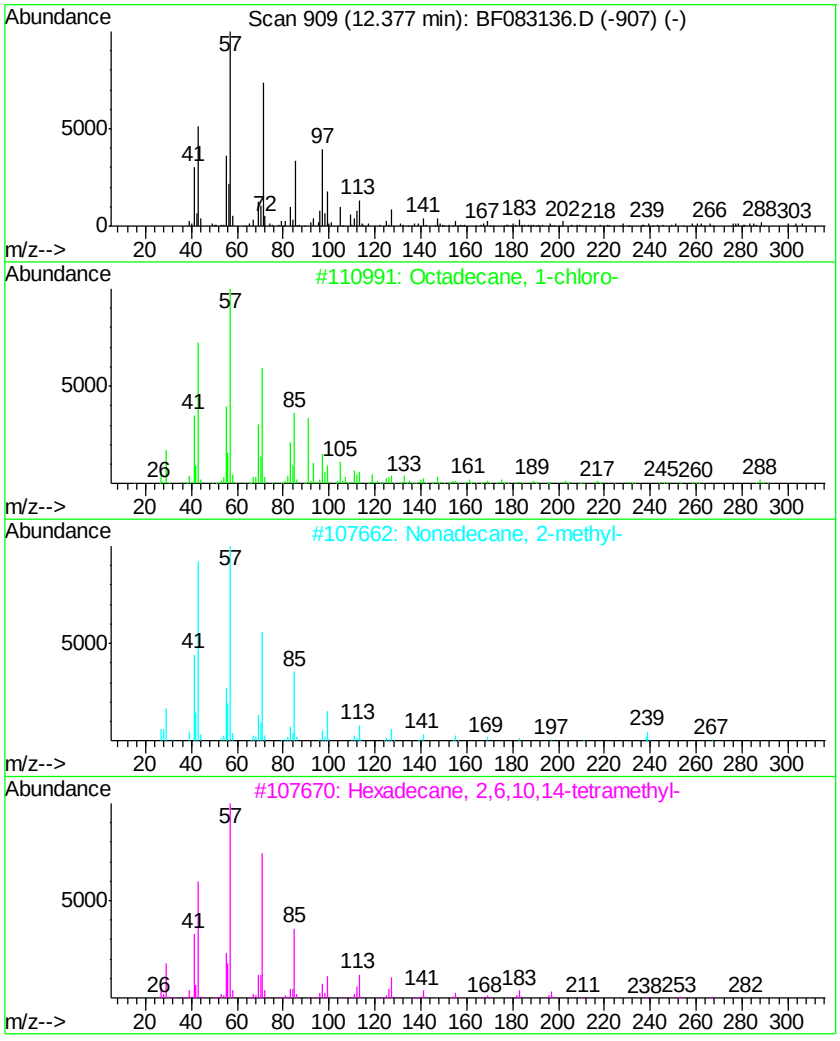
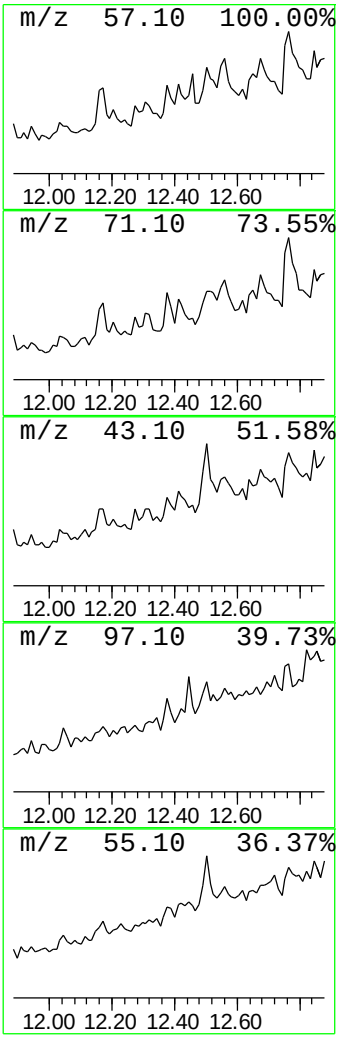
Peak Number 12 Octadecane, 1-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.38	4.05 ng	321618	Phenanthrene-d10	11.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	74
2	Nonadecane, 2-methyl-	282	C20H42	001560-86-7	72
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	68
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	68
5	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	64





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Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
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TIC Library : C:\DATABASE\NIST02.L
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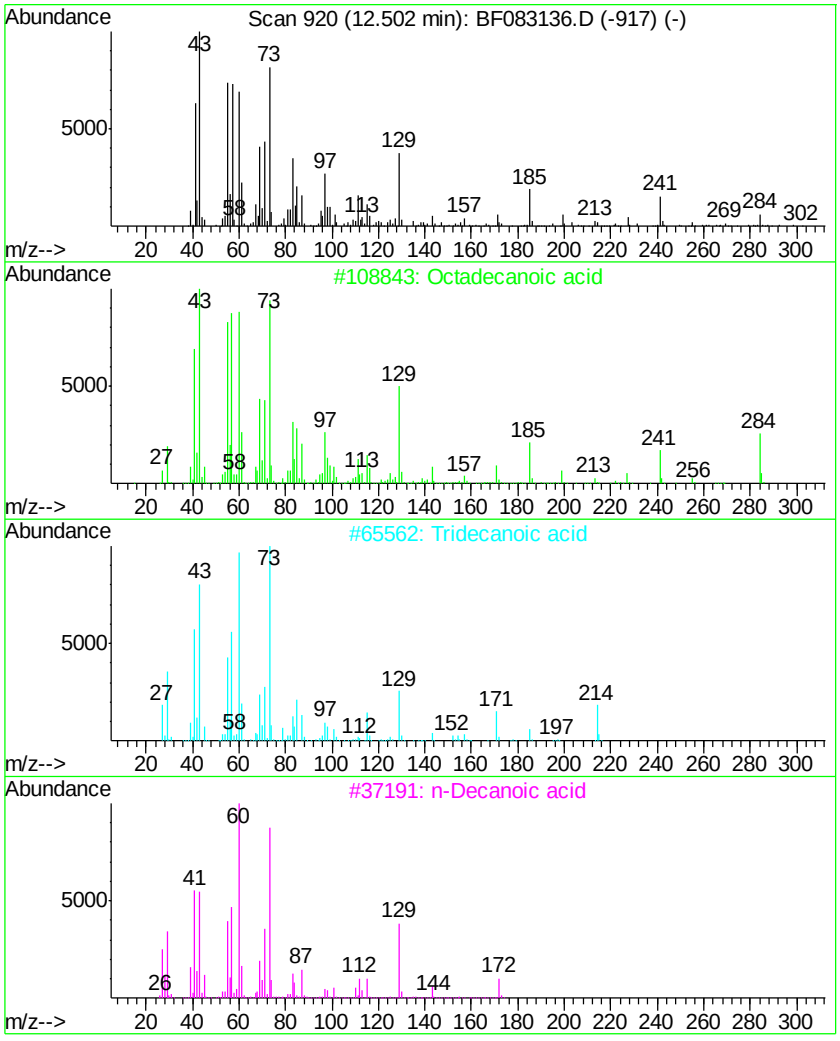
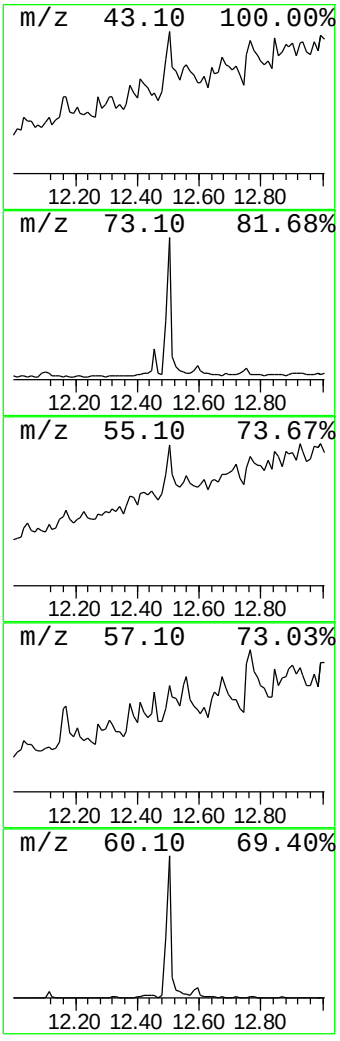
Peak Number 13 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	12.25 ng	852850	Chrysene-d12	13.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	94
2	Tridecanoic acid	214	C13H26O2	000638-53-9	70
3	n-Decanoic acid	172	C10H20O2	000334-48-5	64
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	62
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	46





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Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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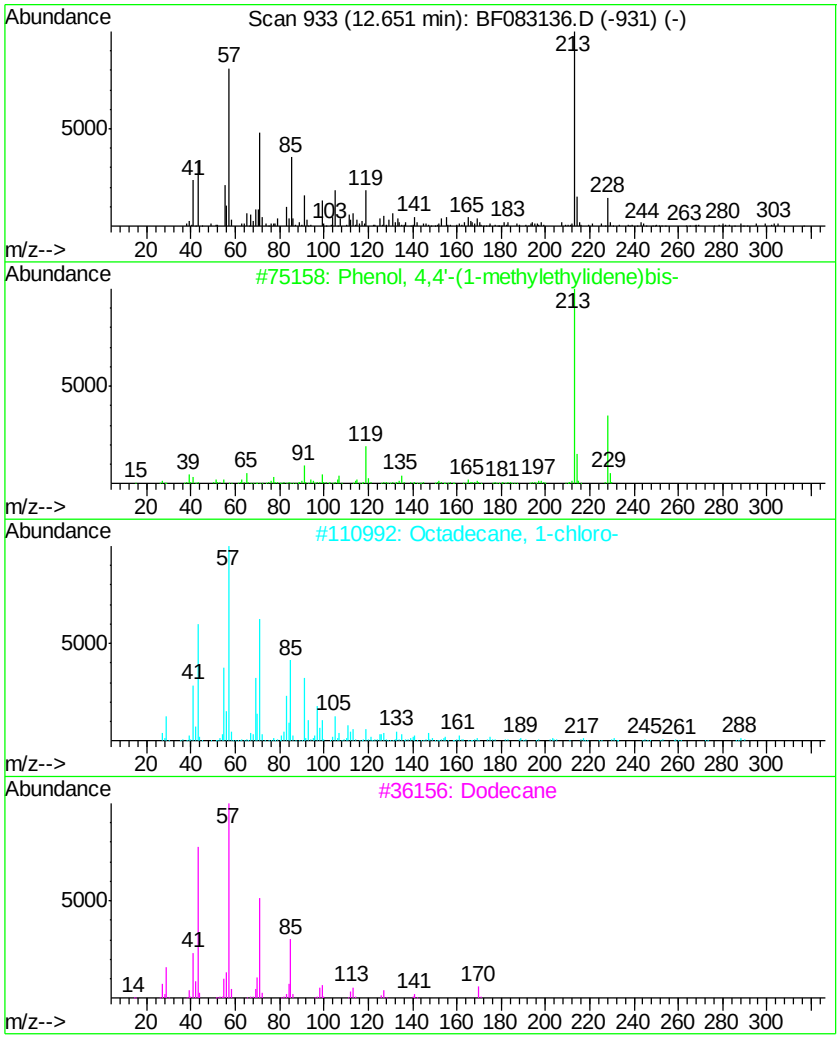
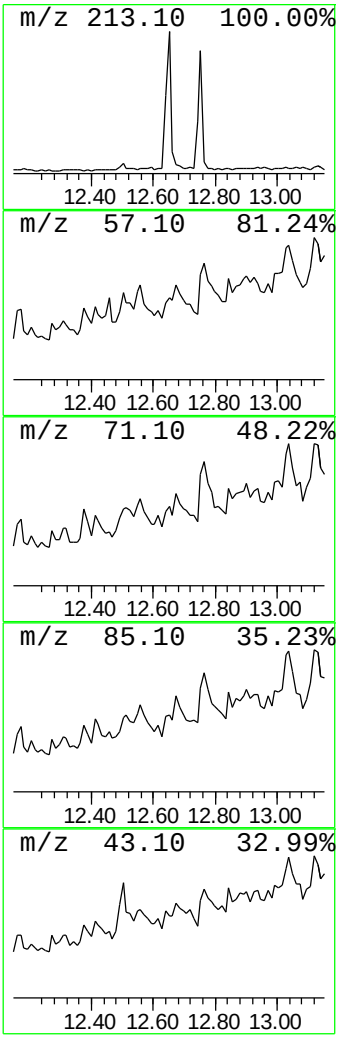
Peak Number 14 unknown12.65 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	4.03 ng	280725	Chrysene-d12	13.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	49
2	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	46
3	Dodecane	170	C12H26	000112-40-3	42
4	Stannane, chlorotriethyl-	242	C6H15ClSn	000994-31-0	42
5	Octacosane	394	C28H58	000630-02-4	38





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Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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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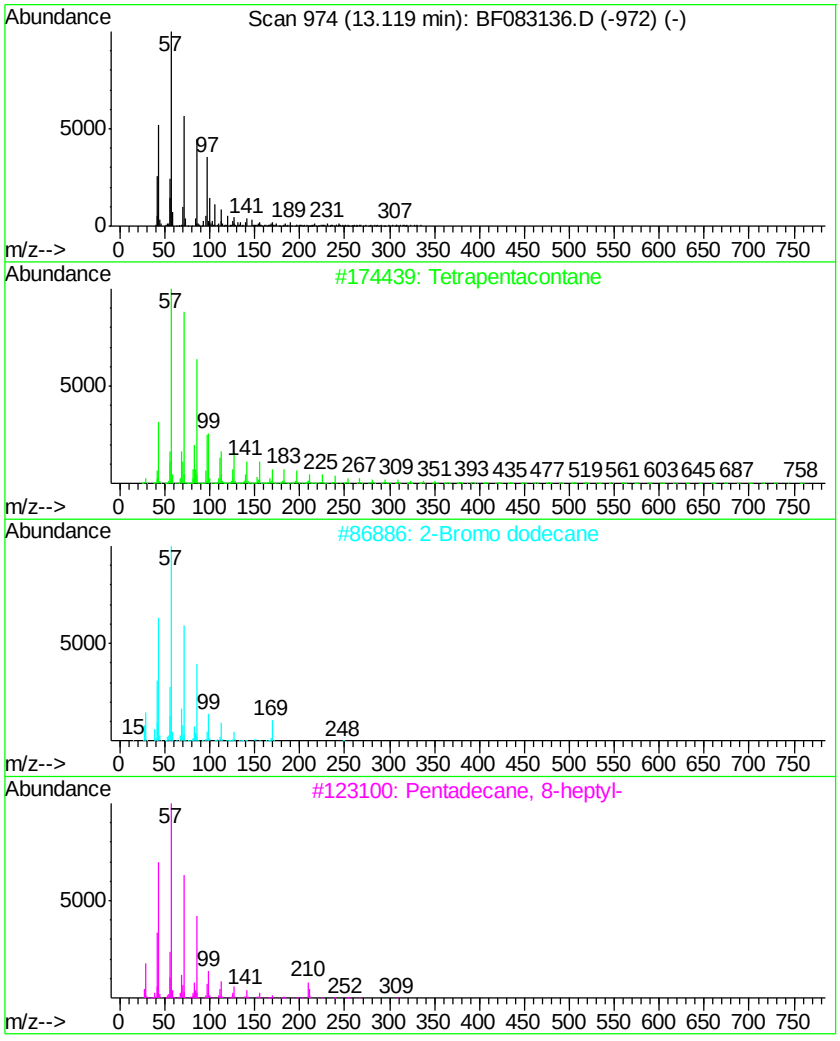
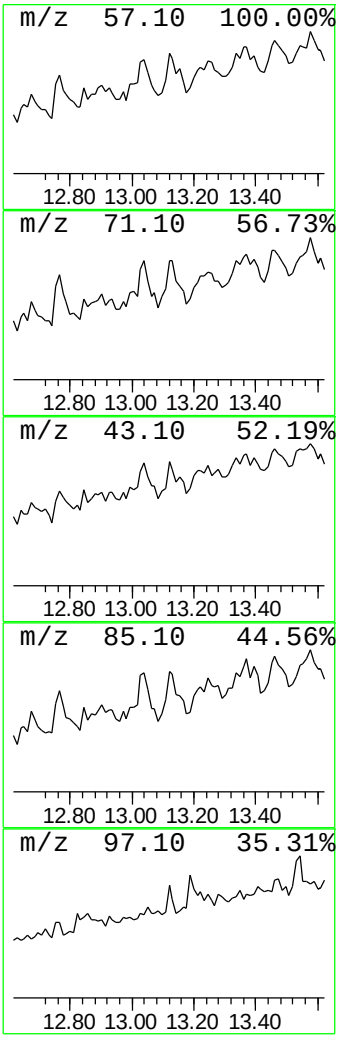
Peak Number 15 Tetrapentacontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	6.07 ng	422680	Chrysene-d12	13.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetrapentacontane	759	C54H110	005856-66-6	72
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	72
3	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	72
4	Heptadecane	240	C17H36	000629-78-7	64
5	Eicosane	282	C20H42	000112-95-8	64





Tentatively Identified Compound (LSC) summary

Data Path : V:\HPCHEM1\BNA_F\DATA\BF112115\
Data File : BF083136.D
Acq On : 22 Nov 2015 4:03
Operator : UM/IZ
Sample : G4468-02
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TEC-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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...	4.81	16.1	ng	957273	1	6.65	1190590	20.0
unknown6.42	6.42	52.2	ng	3109150	1	6.65	1190590	20.0
1-Hexanol, 2-ethyl-	6.75	3.8	ng	227701	1	6.65	1190590	20.0
1-Phenoxypropan-2-ol	8.27	2.0	ng	163329	2	7.94	1626050	20.0
unknown9.37	9.37	12.2	ng	1096580	3	9.69	1795200	20.0
Tridecane	10.60	3.6	ng	284695	4	11.16	1586790	20.0
unknown11.05	11.05	4.9	ng	391955	4	11.16	1586790	20.0
1-Hexadecanethiol	11.66	2.7	ng	216649	4	11.16	1586790	20.0
n-Hexadecanoic acid	11.71	8.7	ng	690985	4	11.16	1586790	20.0
Hexatriacontane	12.17	4.2	ng	330896	4	11.16	1586790	20.0
Octadecane, 1-chl...	12.38	4.0	ng	321618	4	11.16	1586790	20.0
Octadecanoic acid	12.50	12.3	ng	852850	5	13.79	1392800	20.0
unknown12.65	12.65	4.0	ng	280725	5	13.79	1392800	20.0
Tetrapentacontane	13.12	6.1	ng	422680	5	13.79	1392800	20.0