LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

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-BF113015.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	3.138	123	127	134	rVB	138029	288336	4.05%	0.448%
2	4.659	257	260	272	rVB	1136721	2103837	29.58%	3.270%
3	4.933	281	284	286	rBV	691491	993328	13.97%	1.544%
4	5.047	292	294	299	rBV	270028	440274	6.19%	0.684%
5	5.139	299	302	304	rBV	4423996	5451175	76.65%	8.472%
6	5.333	317	319	325	rVB	148873	165912	2.33%	0.258%
7	6.053	380	382	384	rBV	69748	74557	1.05%	0.116%
8	6.213	394	396	399	rBV	4439118	5384828	75.72%	8.369%
9	6.316	403	405	407	rVV	5568051	5322482	74.84%	8.272%
10	6.362	407	409	413	rVB2	93565	182428	2.57%	0.284%
11	6.544	422	425	427	rBV	637450	916502	12.89%	1.424%
12	6.693	436	438	443	rBV	3615575	3972108	55.85%	6.174%
13	6.796	445	447	450	rBV	897821	1046106	14.71%	1.626%
14	7.116	472	475	477	rBV	3441222	3779123	53.14%	5.874%
15	7.607	517	518	525	rVB3	72921	144184	2.03%	0.224%
16	7.847	535	539	542	rBV	4551669	7111861	100.00%	11.053%
17	7.905	542	544	548	rVB	141055	188775	2.65%	0.293%
18	8.545	597	600	602	rBV	223579	257339	3.62%	0.400%
19	8.636	606	608	611	rVB	193288	169739	2.39%	0.264%
20	8.910	629	632	634	rVV	5308139	5457760	76.74%	8.483%
21	9.002	638	640	643	rVV	168915	195219	2.74%	0.303%
22	9.093	645	648	652	rVB	40878	78171	1.10%	0.121%
23	9.310	665	667	669	rVV	1191415	1085142	15.26%	1.687%
24	9.436	674	678	680	rVB	158697	181067	2.55%	0.281%
25	9.528	684	686	688	rVV	186794	211165	2.97%	0.328%
26	9.573	688	690	692	rVV	1551473	1524845	21.44%	2.370%
27	9.608	692	693	695	rVB	158909	119534	1.68%	0.186%
28	9.779	704	708	710	rBV4	43118	115478	1.62%	0.179%
29	9.848	710	714	716	rVV2	114953	200729	2.82%	0.312%
30	9.893	716	718	723	rVB2	79531	124530	1.75%	0.194%
31	10.031	723	730	732	rBV	302458	321716	4.52%	0.500%
32	10.122	736	738	742	rVB	94607	137227	1.93%	0.213%
33	10.248	746	749	752	rBV	92813	159837	2.25%	0.248%
34	10.362	756	759	762	rVB	3245088	3479174	48.92%	5.407%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

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-BF113015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.511	770	772	776	rBV	198311	323641	4.55%	0.503%
36	11.002	813	815	816	rBV2	107661	130652	1.84%	0.203%
37	11.105	822	824	829	rBV2	1094294	1951027	27.43%	3.032%
38	11.196	829	832	838	rVB	144341	204732	2.88%	0.318%
39	11.711	872	877	879	rBV3	46829	101157	1.42%	0.157%
40	11.814	883	886	888	rBV	559223	664732	9.35%	1.033%
41	12.442	938	941	944	rVV2	66852	137524	1.93%	0.214%
42	12.636	956	958	961	rBV	197378	227834	3.20%	0.354%
43	12.774	968	970	974	rVB	59698	86659	1.22%	0.135%
44	12.888	977	980	982	rBV2	208634	301716	4.24%	0.469%
45	12.945	982	985	987	rVV2	309141	465844	6.55%	0.724%
46	12.991	987	989	996	rVB4	29780	104675	1.47%	0.163%
47	13.208	1004	1008	1010	rBV	3284515	4892433	68.79%	7.604%
48	13.437	1026	1028	1033	rVB3	36338	76529	1.08%	0.119%
49	14.637	1130	1133	1137	rBV2	283188	503196	7.08%	0.782%
50	14.728	1137	1141	1143	rVV	980804	1465350	20.60%	2.277%
51	16.214	1268	1271	1274	rVB	135467	195875	2.75%	0.304%
52	16.283	1274	1277	1285	rBV2	30657	90543	1.27%	0.141%
53	16.785	1318	1321	1327	rVB	635297	956395	13.45%	1.486%
54	17.506	1382	1384	1387	rVB4	40063	75967	1.07%	0.118%

Sum of corrected areas: 64340969

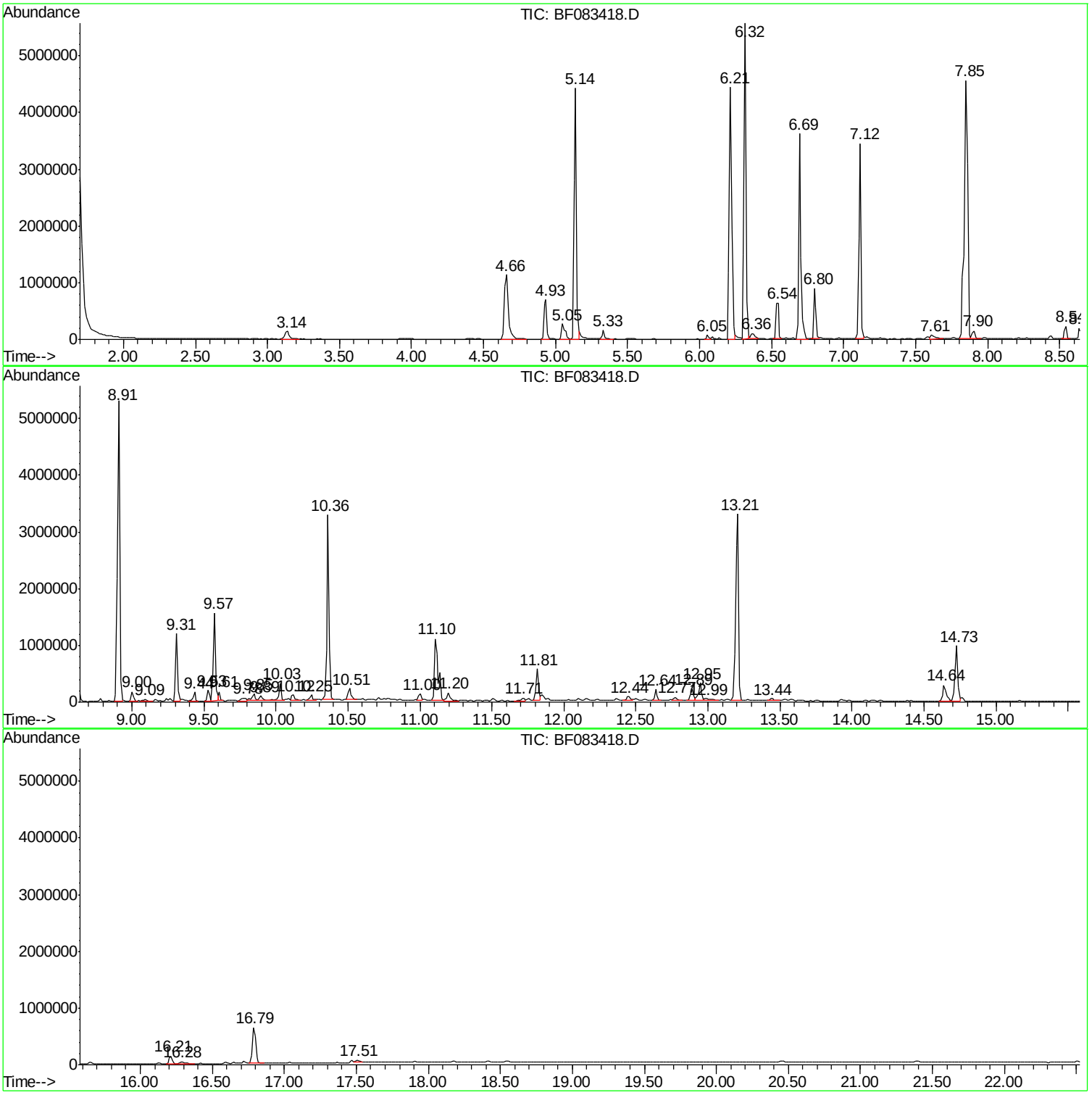
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

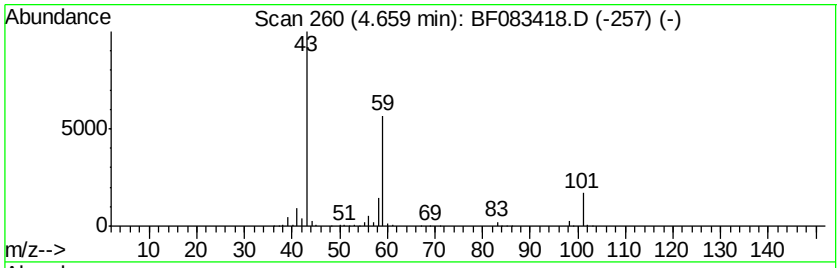
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

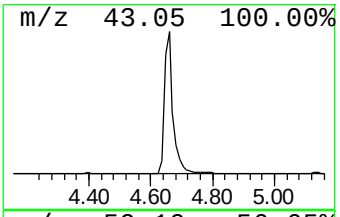
****

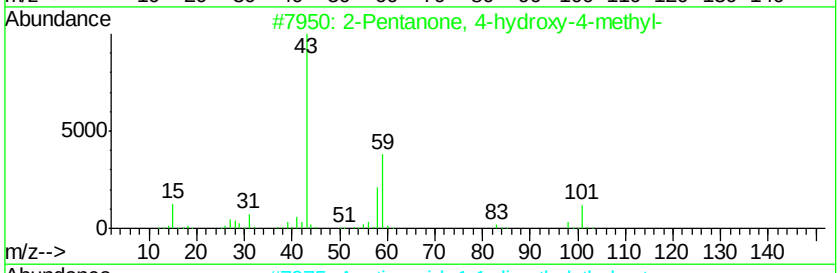
Peak Number 2 unknown4.66 Concentration Rank 3

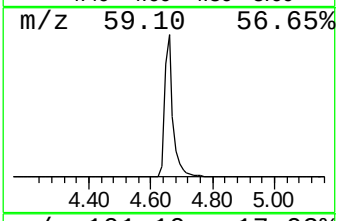
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	45.91 ng	2103840	1,4-Dichlorobenzene-d4	6.54

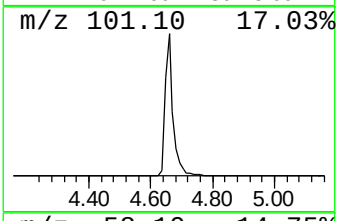
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	1,2-Butanediol, (.+/-.)-	90	C4H10O2	026171-83-5	9

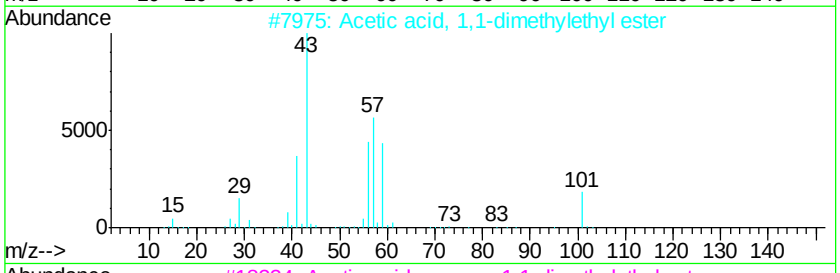


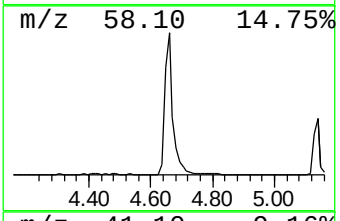


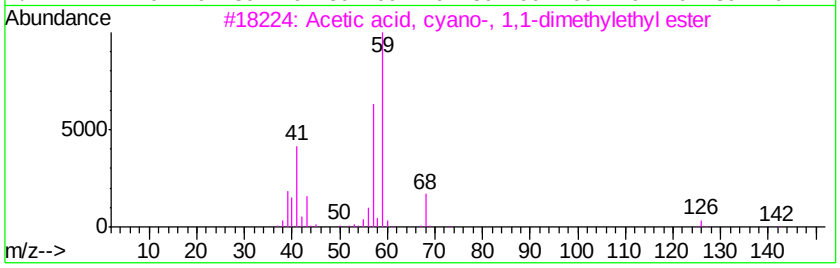


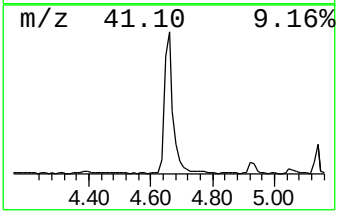












Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

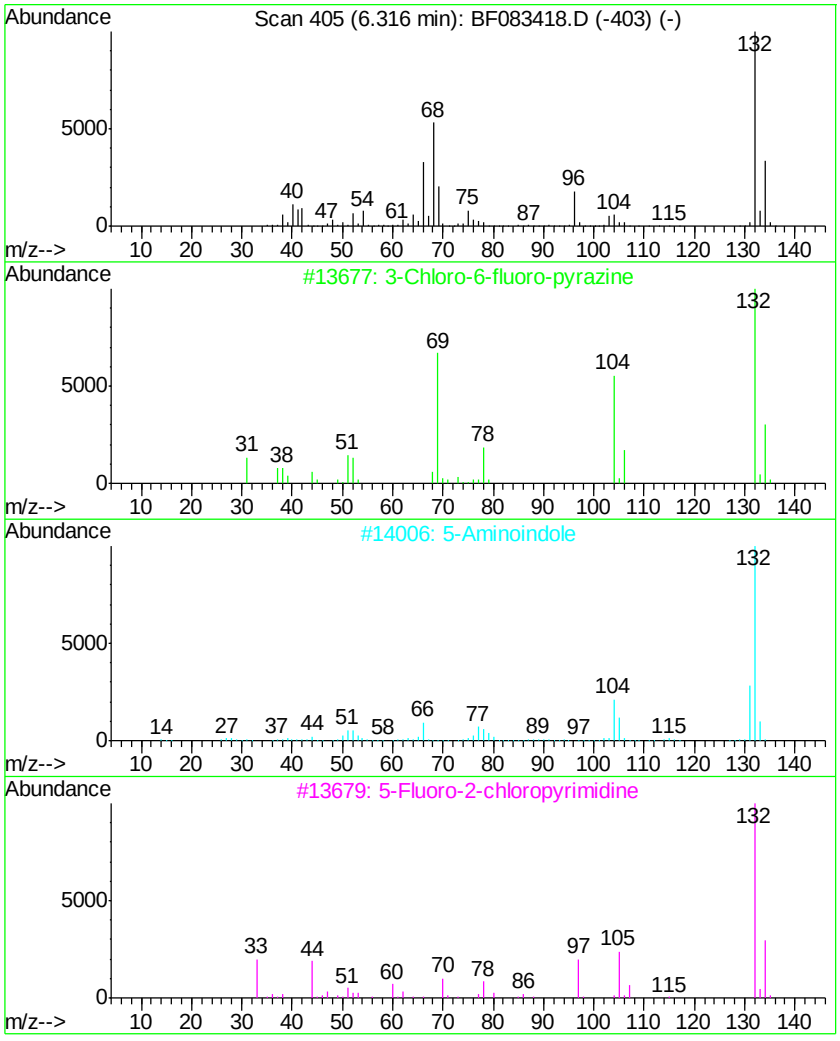
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

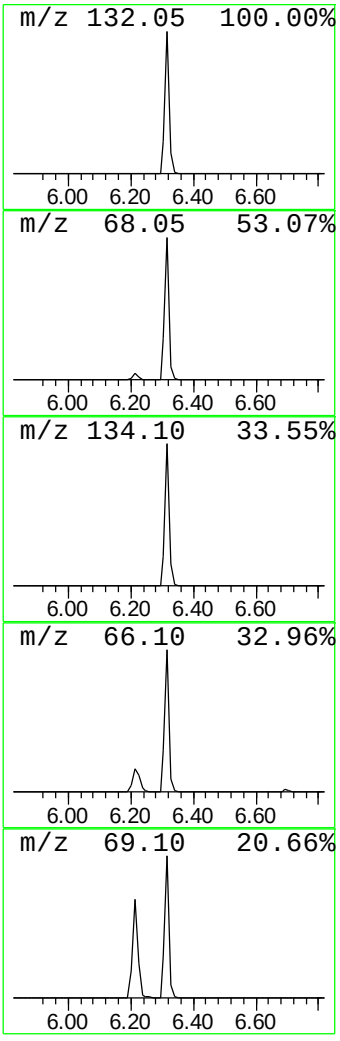
****

Peak Number 6 unknown6.32 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	116.15 ng	5322480	1,4-Dichlorobenzene-d4	6.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole	132	C8H8N2	005192-03-0	22
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
5	(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10





Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

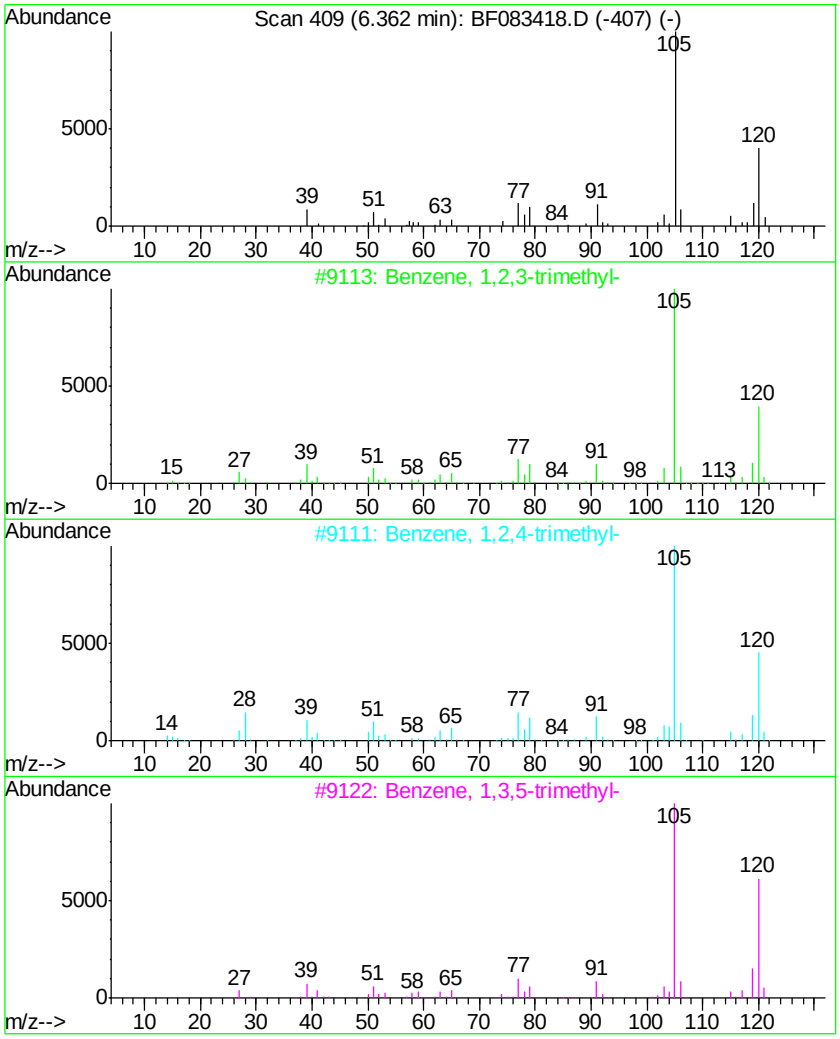
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

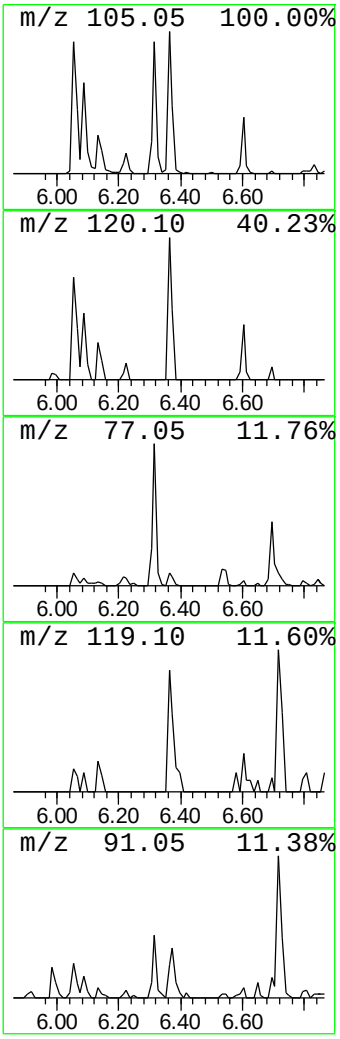
****

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.36	3.98 ng	182428	1,4-Dichlorobenzene-d4	6.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

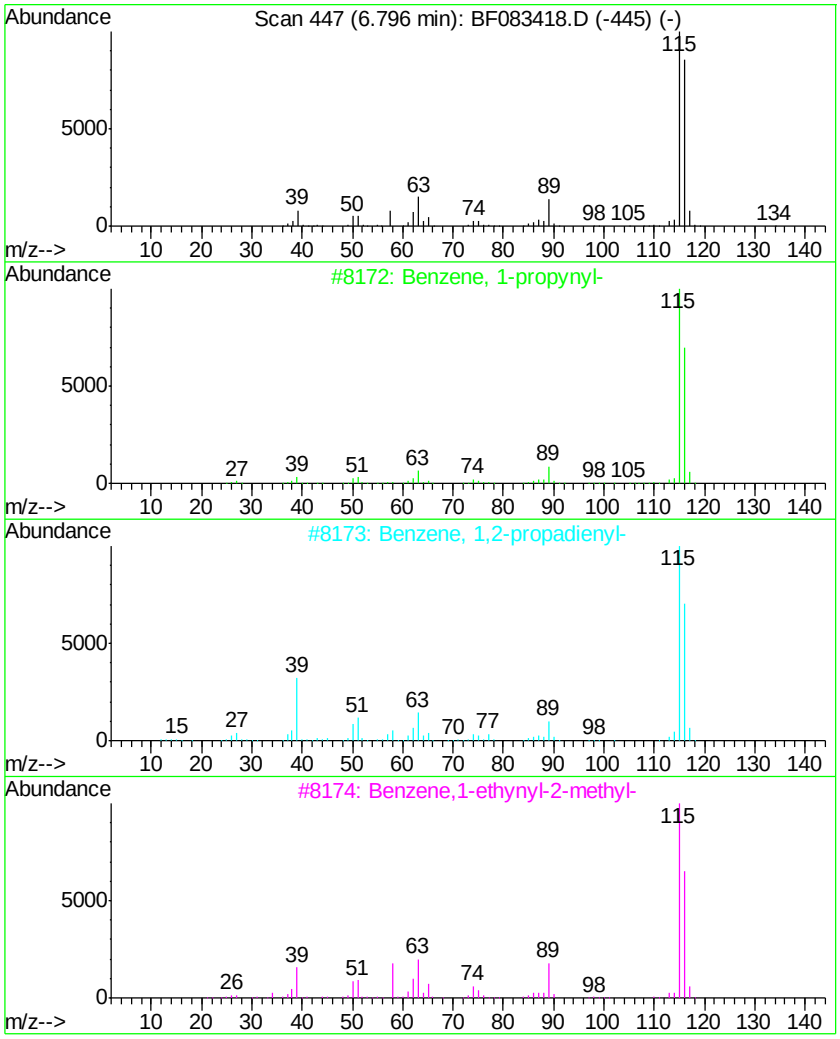
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

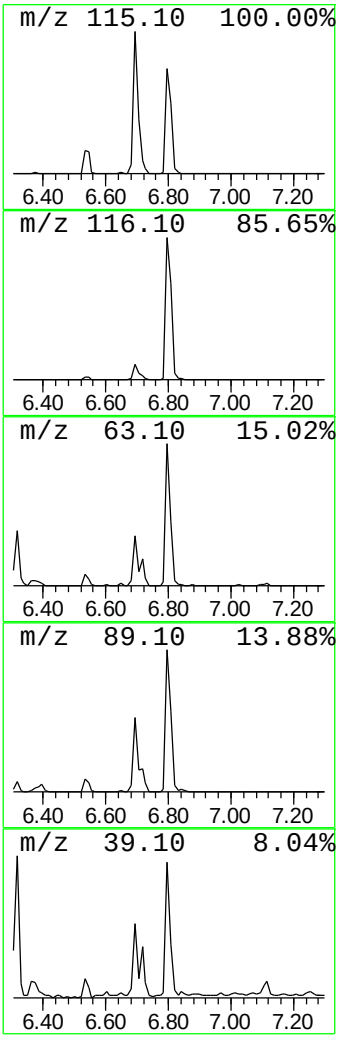
****

Peak Number 9 Benzene, 1-propynyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	22.83 ng	1046110	1,4-Dichlorobenzene-d4	6.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
2	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
3	Benzene,1-ethynyl-2-methyl-	116	C9H8	000766-47-2	91
4	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
5	Indene	116	C9H8	000095-13-6	81





Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

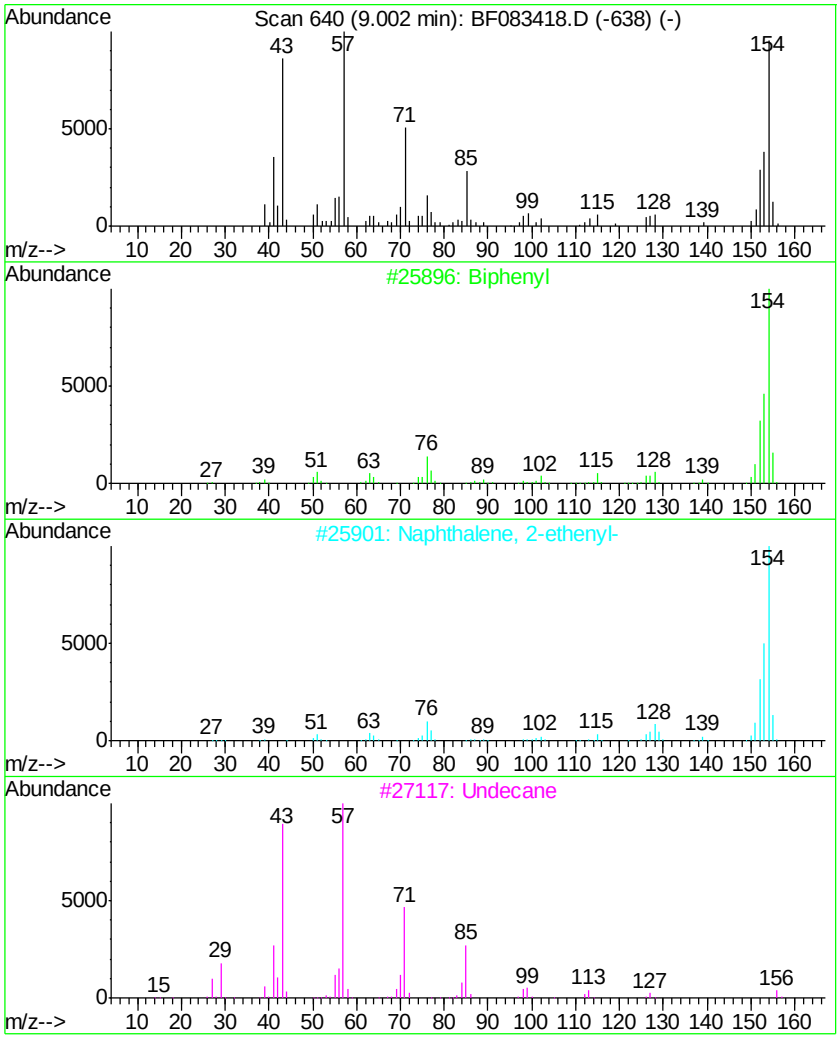
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

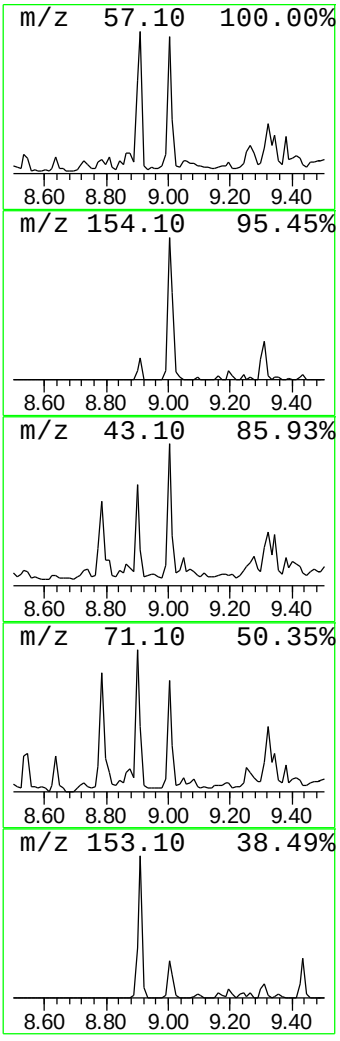
****

Peak Number 10 Biphenyl Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.00	2.56 ng	195219	Acenaphthene-d10	9.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenyl	154	C12H10	000092-52-4	91
2	Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	55
3	Undecane	156	C11H24	001120-21-4	35
4	Propanedinitrile, (phenylmethyle...	154	C10H6N2	002700-22-3	30
5	Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	25





Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

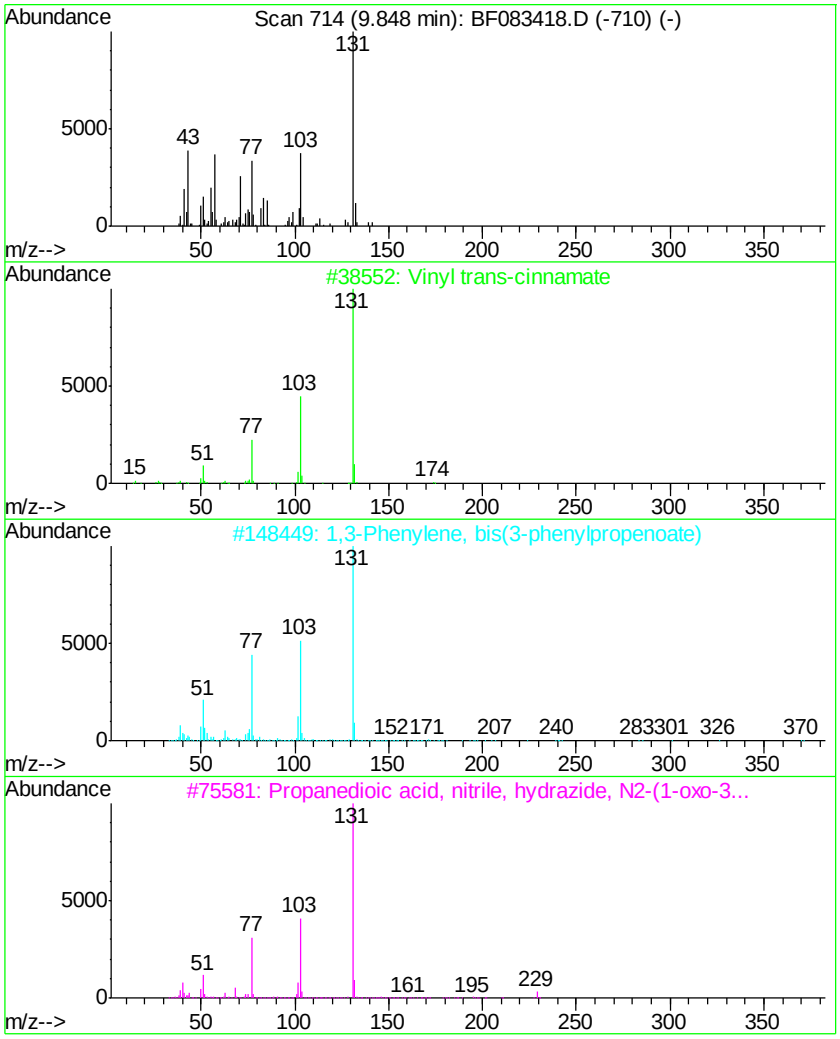
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

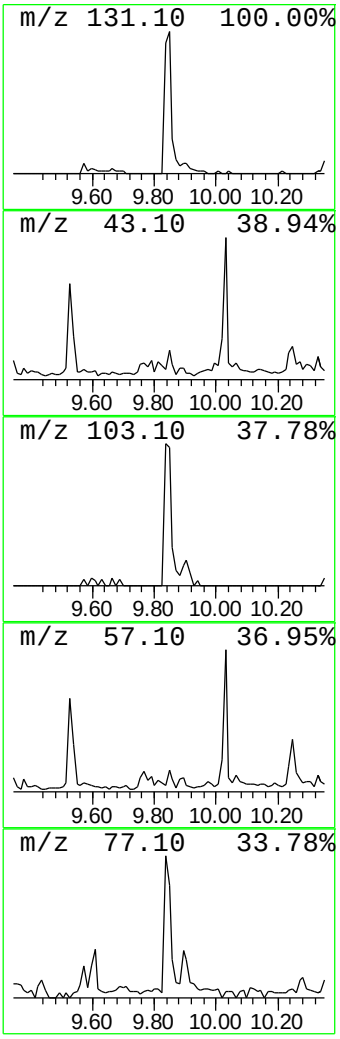
****

Peak Number 11 Vinyl trans-cinnamate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	2.63 ng	200729	Acenaphthene-d10	9.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vinyl trans-cinnamate	174	C11H10O2	017719-70-9	59
2	1,3-Phenylene, bis(3-phenylprope...	370	C24H18O4	1000259-62-4	59
3	Propanedioic acid, nitrile, hydr...	229	C12H11N3O2	294878-35-6	45
4	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	42
5	Oxalic acid, monoamide monohydra...	323	C18H17N3O3	1000294-01-4	42





Data Path : Z:\HPCHEM1\BNA F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

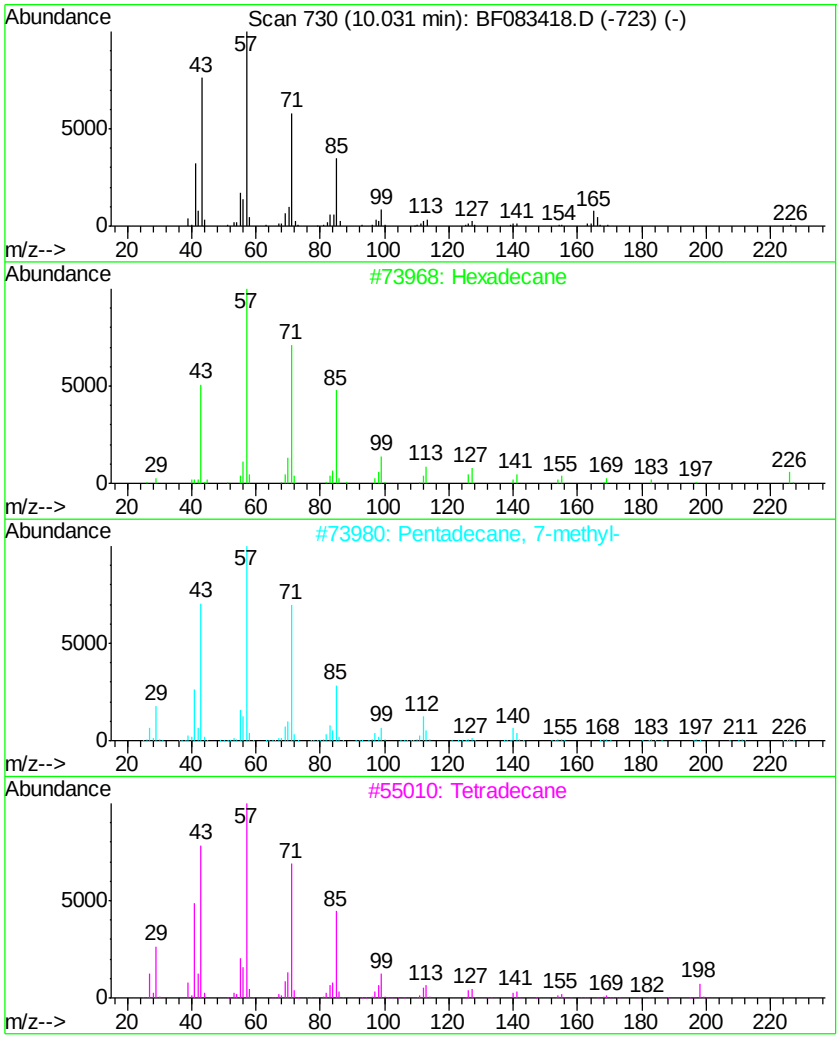
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

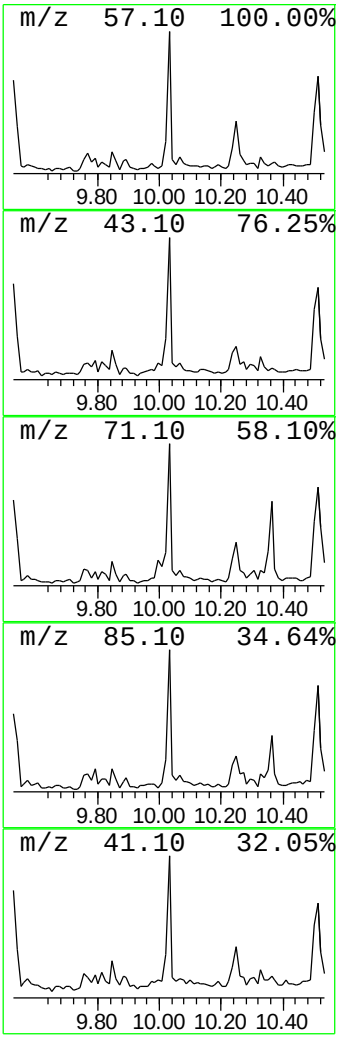
****

Peak Number 12 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.03	4.22 ng	321716	Acenaphthene-d10	9.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	90
3	Tetradecane	198	C14H30	000629-59-4	78
4	Nonane, 5-butyl-	184	C13H28	017312-63-9	64
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	64





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Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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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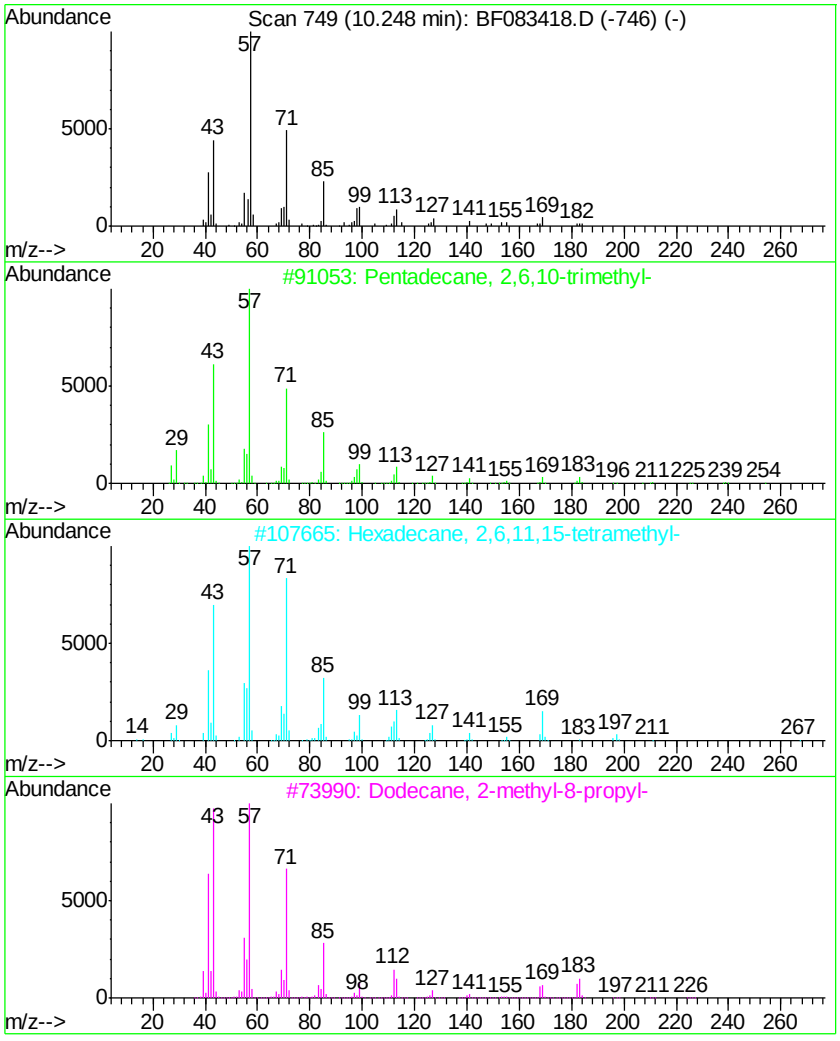
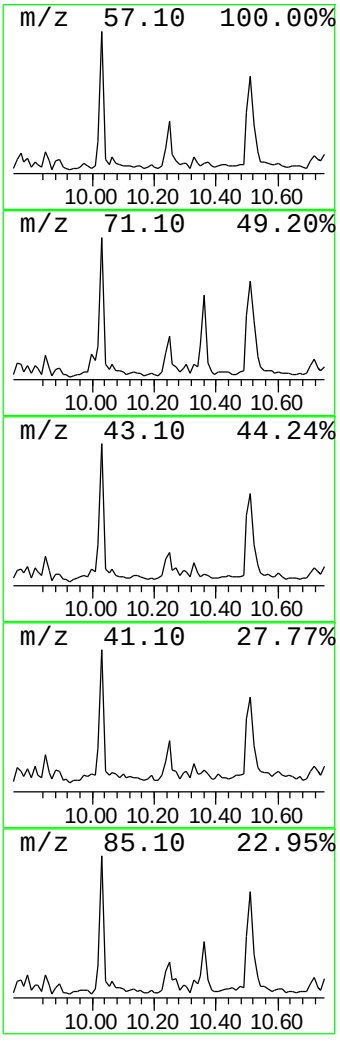
Peak Number 13 Pentadecane, 2,6,10-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	2.10 ng	159837	Acenaphthene-d10	9.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	91
2	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	68
3	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	64
4	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	64
5	Octacosane	394	C28H58	000630-02-4	64





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Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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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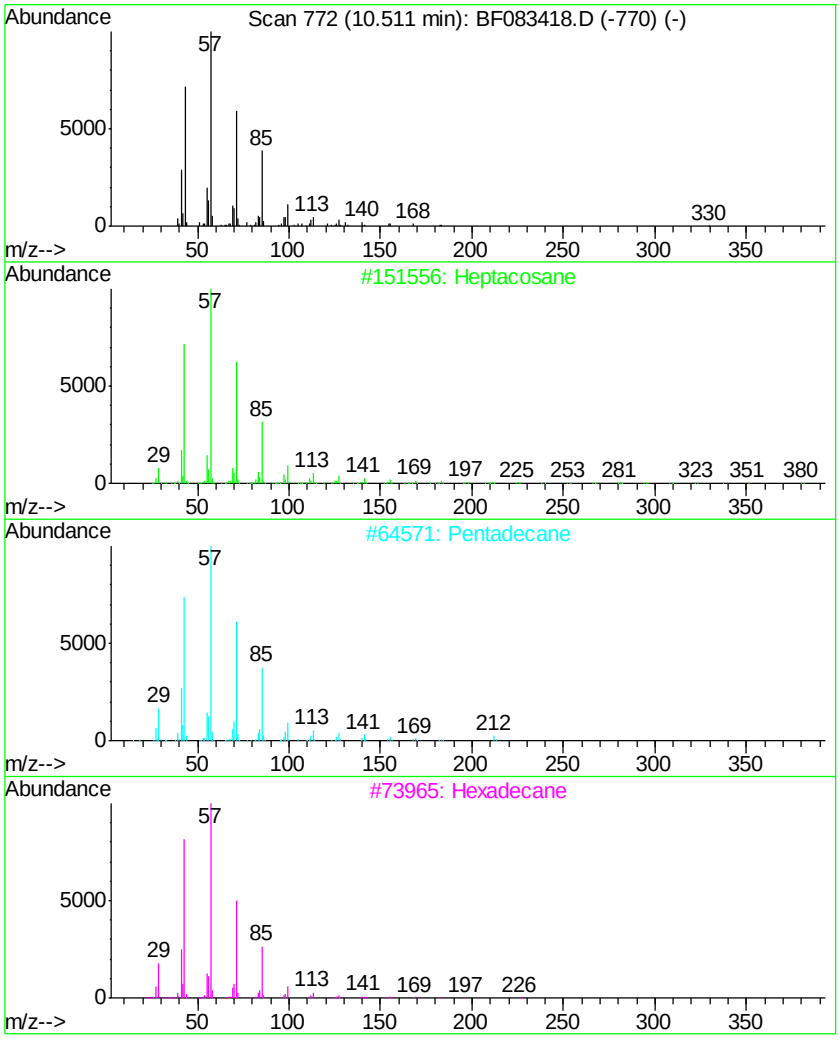
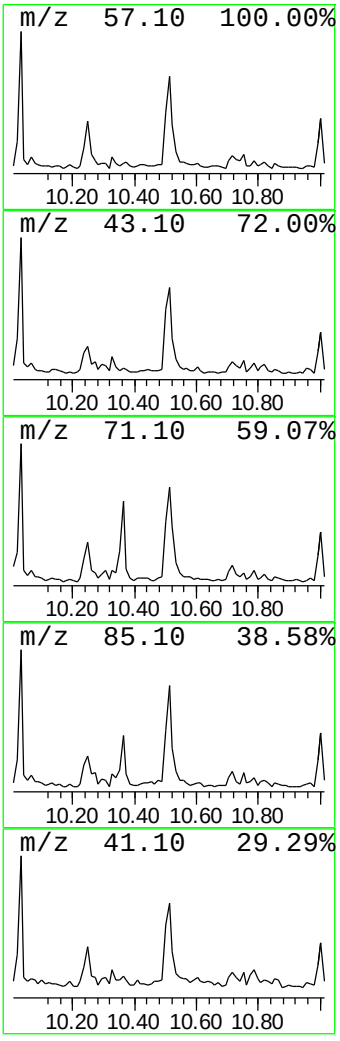
Peak Number 14 Heptacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	3.32 ng	323641	Phenanthrene-d10	11.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	90
2	Pentadecane	212	C15H32	000629-62-9	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tetracosane	338	C24H50	000646-31-1	90
5	Tetratetracontane	619	C44H90	007098-22-8	90





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Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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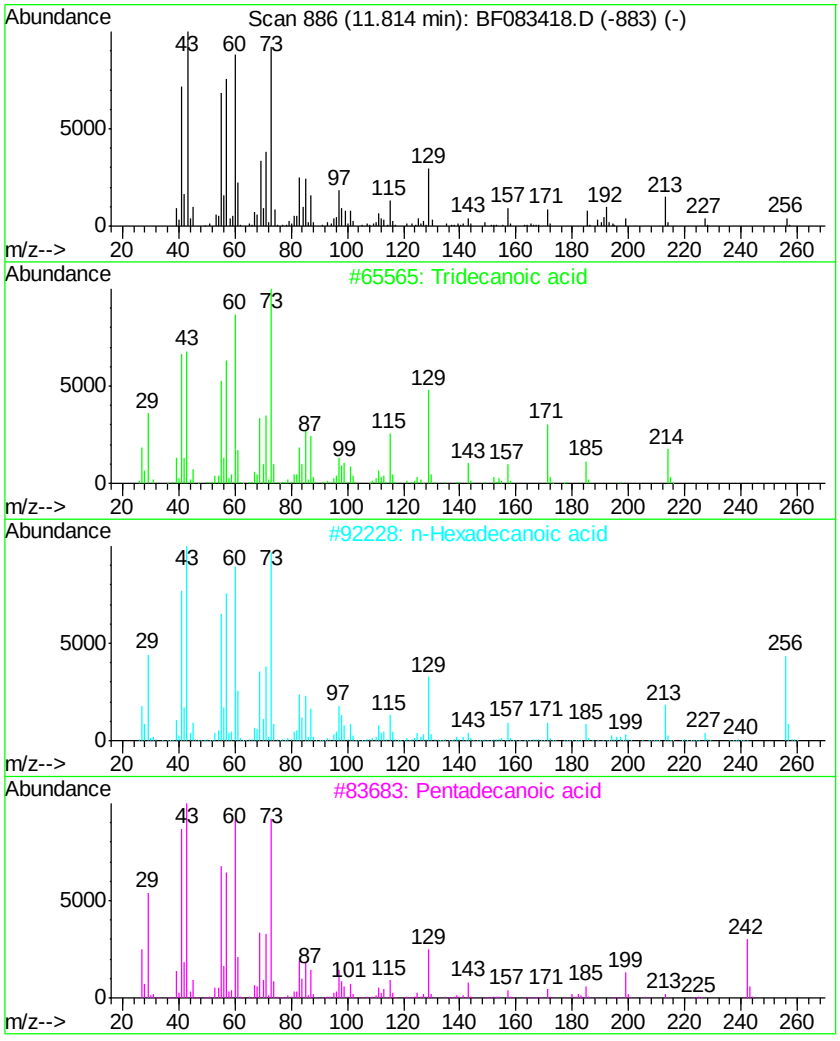
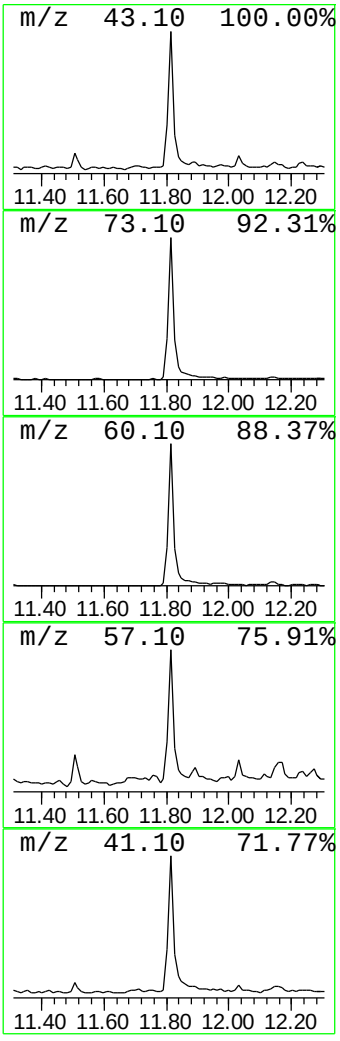
Peak Number 16 Tridecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	6.81 ng	664732	Phenanthrene-d10	11.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	93
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	35
5	Undecane	156	C11H24	001120-21-4	18





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Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.M
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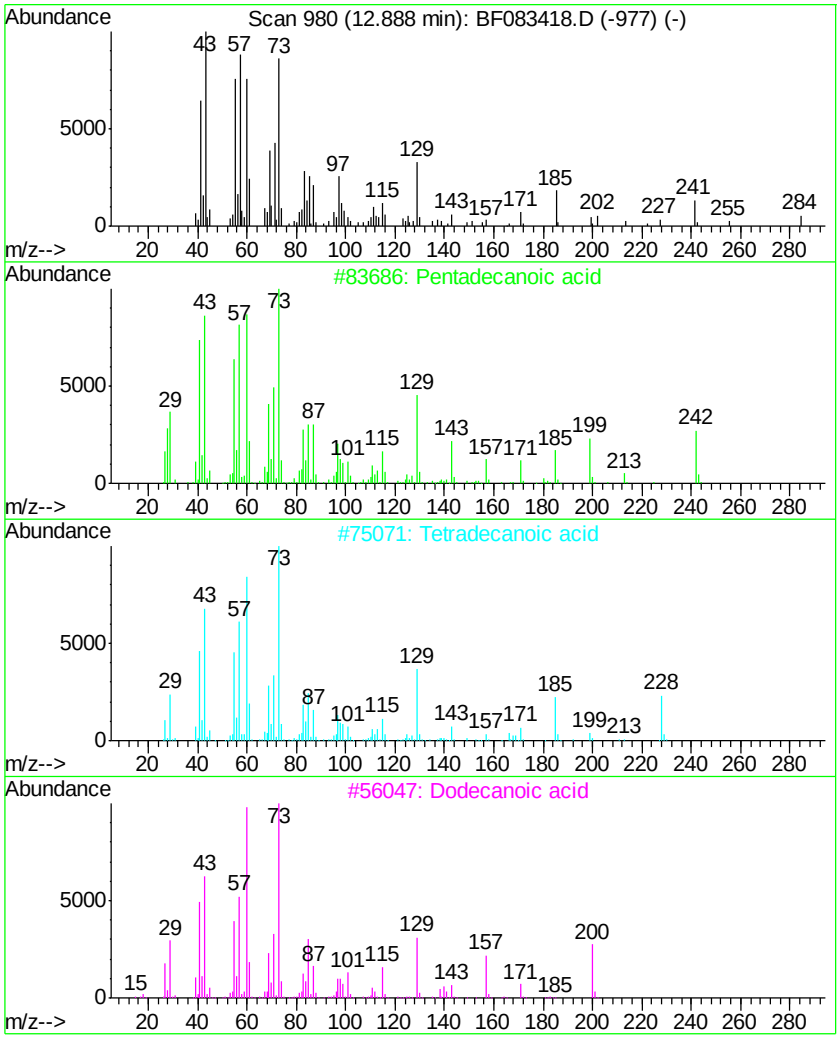
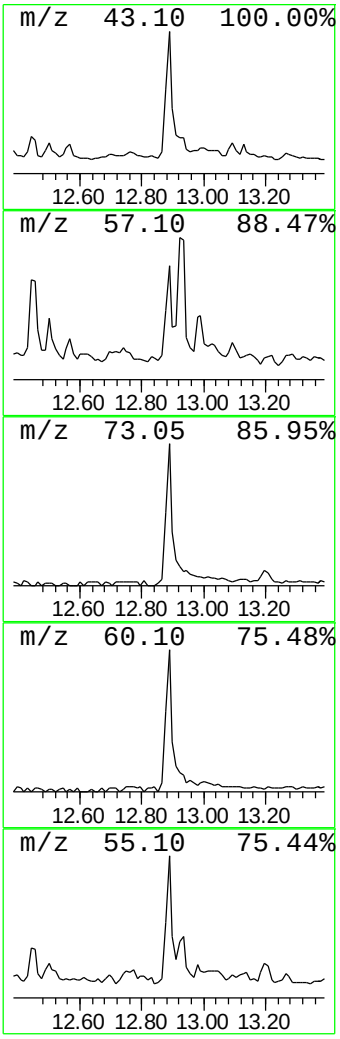
Peak Number 17 Pentadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	3.09 ng	301716	Phenanthrene-d10	11.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3	Dodecanoic acid	200	C12H24O2	000143-07-7	59
4	1-Eicosanol	298	C20H42O	000629-96-9	14
5	Dodecane	170	C12H26	000112-40-3	11





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Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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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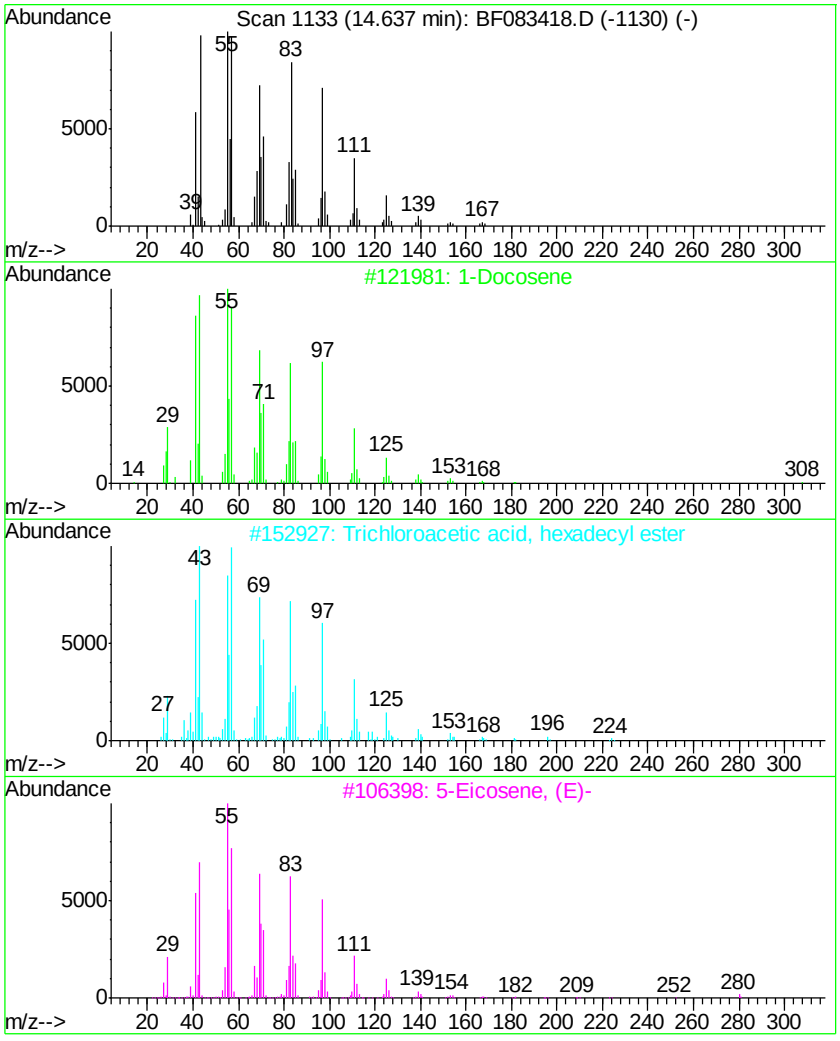
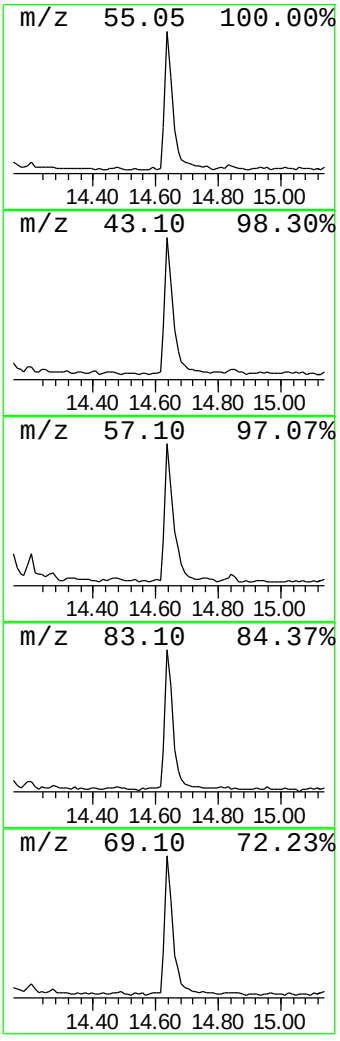
Peak Number 18 1-Docosene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	6.87 ng	503196	Chrysene-d12	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3	5-Eicosene, (E)-	280	C20H40	074685-30-6	91
4	1-Nonadecene	266	C19H38	018435-45-5	91
5	2- Bromopropionic acid, pentadec...	362	C18H35BrO2	1000292-44-2	91





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Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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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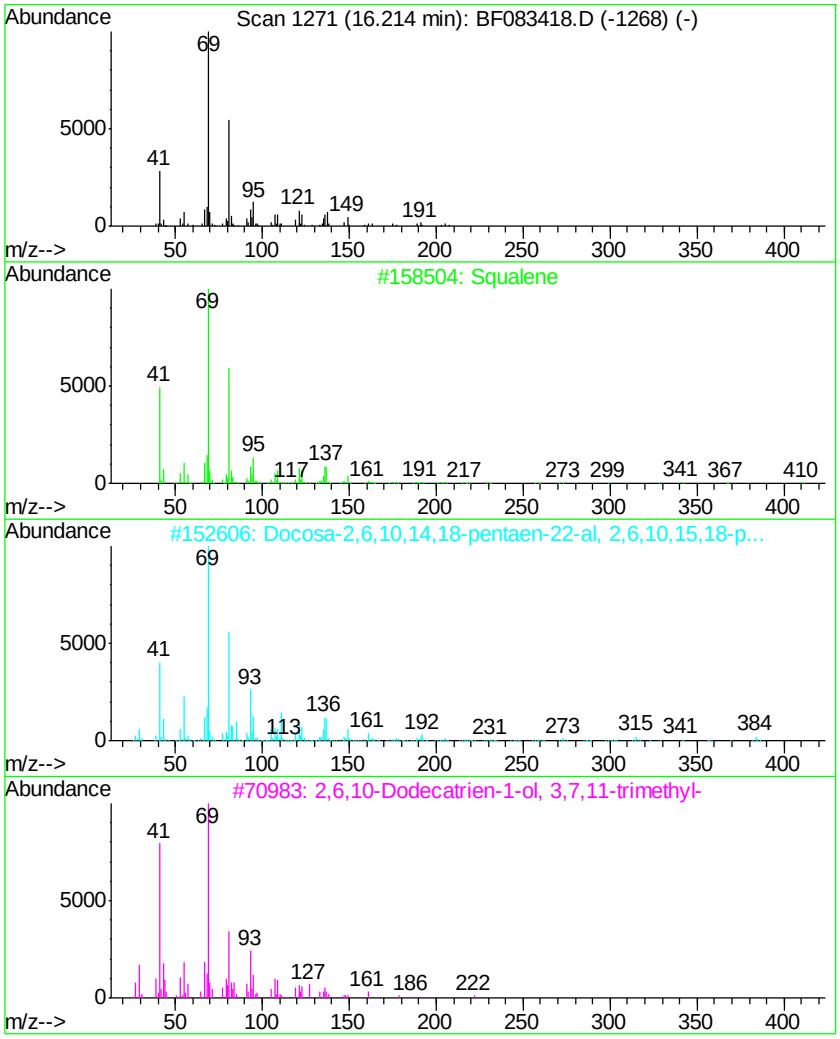
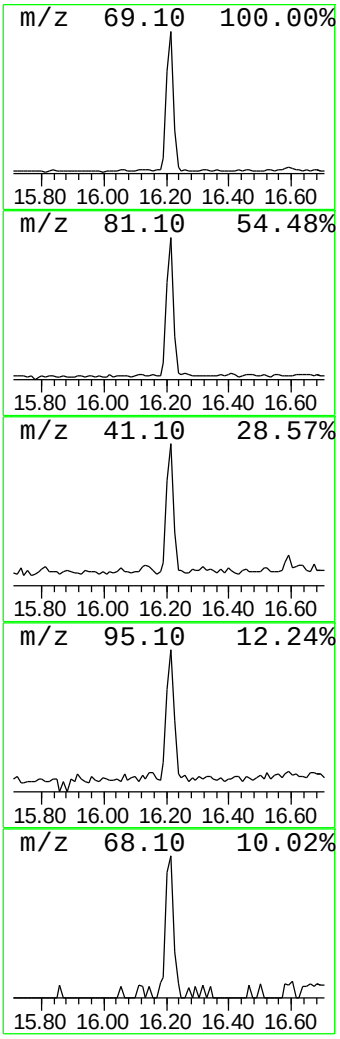
Peak Number 19 Squalene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.21	4.10 ng	195875	Perylene-d12	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	007683-64-9	91
2	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	83
3	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	78
4	4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	56
5	1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120215\
Data File : BF083418.D
Acq On : 2 Dec 2015 15:11
Operator : UM/IZ
Sample : G4582-16
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-P3WC-09(8-14)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF113015.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
unknown4.66	4.66	45.9	ng	2103840	1	6.54	916502	20.0
unknown6.32	6.32	116.2	ng	5322480	1	6.54	916502	20.0
Benzene, 1,2,3-tr...	6.36	4.0	ng	182428	1	6.54	916502	20.0
Benzene, 1-propynyl-	6.80	22.8	ng	1046110	1	6.54	916502	20.0
Biphenyl	9.00	2.6	ng	195219	3	9.57	1524850	20.0
Vinyl trans-cinna...	9.85	2.6	ng	200729	3	9.57	1524850	20.0
Hexadecane	10.03	4.2	ng	321716	3	9.57	1524850	20.0
Pentadecane, 2,6,...	10.25	2.1	ng	159837	3	9.57	1524850	20.0
Heptacosane	10.51	3.3	ng	323641	4	11.10	1951030	20.0
Tridecanoic acid	11.81	6.8	ng	664732	4	11.10	1951030	20.0
Pentadecanoic acid	12.89	3.1	ng	301716	4	11.10	1951030	20.0
1-Docosene	14.64	6.9	ng	503196	5	14.73	1465350	20.0
Squalene	16.21	4.1	ng	195875	6	16.79	956395	20.0