LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

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-BF060716.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	1.724	10	12	21	rVB	374033	668632	23.11%	1.987%
2	1.850	21	23	36	rVB	1580261	2893123	100.00%	8.599%
3	3.050	126	128	145	rBV	62586	133432	4.61%	0.397%
4	4.959	293	295	301	rBV	130427	181484	6.27%	0.539%
5	5.381	329	332	334	rBV	2322264	2611496	90.27%	7.762%
6	6.422	420	423	426	rBV	1942546	2415339	83.49%	7.179%
7	6.547	431	434	437	rBV	2200900	2503952	86.55%	7.443%
8	6.776	452	454	457	rBV	631711	627514	21.69%	1.865%
9	6.936	465	468	470	rBV	2173744	1922021	66.43%	5.713%
10	7.199	487	491	493	rBV	33422	41557	1.44%	0.124%
11	7.347	501	504	506	rBV	1579626	1665981	57.58%	4.952%
12	7.450	509	513	520	rBV	32068	47838	1.65%	0.142%
13	7.805	541	544	553	rVB2	33810	52546	1.82%	0.156%
14	8.067	564	567	569	rBV	1017775	927951	32.07%	2.758%
15	9.153	659	662	664	rVV	2287524	2826387	97.69%	8.401%
16	9.233	665	669	670	rVV2	36207	52870	1.83%	0.157%
17	9.268	670	672	677	rVB2	49646	71296	2.46%	0.212%
18	9.485	687	691	694	rBV2	13190	29390	1.02%	0.087%
19	9.542	694	696	698	rVV	935989	1205258	41.66%	3.582%
20	9.759	714	715	718	rVB	36245	30274	1.05%	0.090%
21	9.828	718	721	722	rVV	704390	990682	34.24%	2.945%
22	9.850	722	723	725	rVB	41484	34152	1.18%	0.102%
23	10.239	754	757	758	rBV	29885	41901	1.45%	0.125%
24	10.262	758	759	760	rVB	52042	35682	1.23%	0.106%
25	10.365	767	768	772	rVB2	51908	45480	1.57%	0.135%
26	10.479	775	778	781	rBV	31253	45538	1.57%	0.135%
27	10.616	786	790	792	rBV	1335976	1806758	62.45%	5.370%
28	10.731	799	800	805	rBV	69654	103122	3.56%	0.307%
29	11.176	835	839	841	rBV2	51451	97436	3.37%	0.290%
30	11.211	841	842	844	rVB	54446	45272	1.56%	0.135%
31	11.302	848	850	854	rVV2	1112040	1204438	41.63%	3.580%
32	11.371	854	856	858	rVB2	39572	54697	1.89%	0.163%
33	11.542	868	871	874	rBV4	57284	149343	5.16%	0.444%
34	11.611	874	877	878	rVV	58513	70541	2.44%	0.210%

LSC Area Percent Report

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Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
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Start Thrs: 0.2 Max Peaks: 100
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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.805	892	894	895	rBV	33823	38437	1.33%	0.114%
36	11.839	895	897	901	rVV3	64377	157924	5.46%	0.469%
37	11.931	901	905	908	rVV3	42099	112557	3.89%	0.335%
38	12.022	911	913	914	rBV	61139	90531	3.13%	0.269%
39	12.091	917	919	921	rVV3	34084	63573	2.20%	0.189%
40	12.251	931	933	935	rVB3	39333	39841	1.38%	0.118%
41	12.399	944	946	949	rVB2	71723	89736	3.10%	0.267%
42	12.514	954	956	958	rBV	274738	284119	9.82%	0.845%
43	12.559	958	960	962	rVB	143703	123519	4.27%	0.367%
44	12.719	972	974	977	rVB2	396702	485548	16.78%	1.443%
45	12.777	977	979	983	rVB3	77406	148629	5.14%	0.442%
46	12.891	987	989	992	rVB	2120185	2090846	72.27%	6.215%
47	12.959	992	995	996	rBV2	24661	36175	1.25%	0.108%
48	13.028	999	1001	1003	rBV	58399	75951	2.63%	0.226%
49	13.131	1008	1010	1011	rBV	111572	103050	3.56%	0.306%
50	13.165	1011	1013	1018	rVB5	40360	70125	2.42%	0.208%
51	13.428	1034	1036	1037	rBV	197463	215445	7.45%	0.640%
52	13.474	1037	1040	1041	rVV	97015	114358	3.95%	0.340%
53	13.794	1066	1068	1071	rVB2	163756	206108	7.12%	0.613%
54	13.885	1074	1076	1078	rBV	120688	108719	3.76%	0.323%
55	13.931	1078	1080	1085	rVB2	793191	1049331	36.27%	3.119%
56	14.114	1094	1096	1098	rVB	106155	99644	3.44%	0.296%
57	14.422	1121	1123	1124	rBV	59782	83265	2.88%	0.247%
58	14.708	1146	1148	1151	rVB	45506	69419	2.40%	0.206%
59	14.948	1166	1169	1172	rVV	80040	164577	5.69%	0.489%
60	15.028	1174	1176	1179	rVB2	56844	79324	2.74%	0.236%
61	15.222	1191	1193	1196	rVV	49983	70367	2.43%	0.209%
62	15.280	1196	1198	1201	rVV	61587	94588	3.27%	0.281%
63	15.348	1201	1204	1206	rVV	450043	711161	24.58%	2.114%
64	15.382	1206	1207	1211	rVV2	51283	57344	1.98%	0.170%
65	15.474	1213	1215	1220	rVB4	23650	41262	1.43%	0.123%
66	15.668	1227	1232	1240	rVB2	62730	270845	9.36%	0.805%
67	15.783	1240	1242	1244	rBV	23100	32326	1.12%	0.096%
68	15.988	1258	1260	1264	rVB2	25307	46556	1.61%	0.138%
69	16.251	1280	1283	1285	rBV2	16321	32276	1.12%	0.096%
70	16.388	1293	1295	1302	rVB3	30984	73983	2.56%	0.220%
71	16.685	1318	1321	1325	rVB2	28298	54293	1.88%	0.161%

LSC Area Percent Report

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Misc :
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Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

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

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	17.097	1354	1357	1362	rVB	23809	49257	1.70%	0.146%
73	17.657	1403	1406	1410	rBV5	13221	44079	1.52%	0.131%
74	17.748	1411	1414	1419	rVB3	21253	59300	2.05%	0.176%
75	18.011	1430	1437	1448	rVB7	24315	134425	4.65%	0.400%
76	18.423	1470	1473	1480	rVB7	13395	36938	1.28%	0.110%

Sum of corrected areas: 33643134

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Sample : H3437-05
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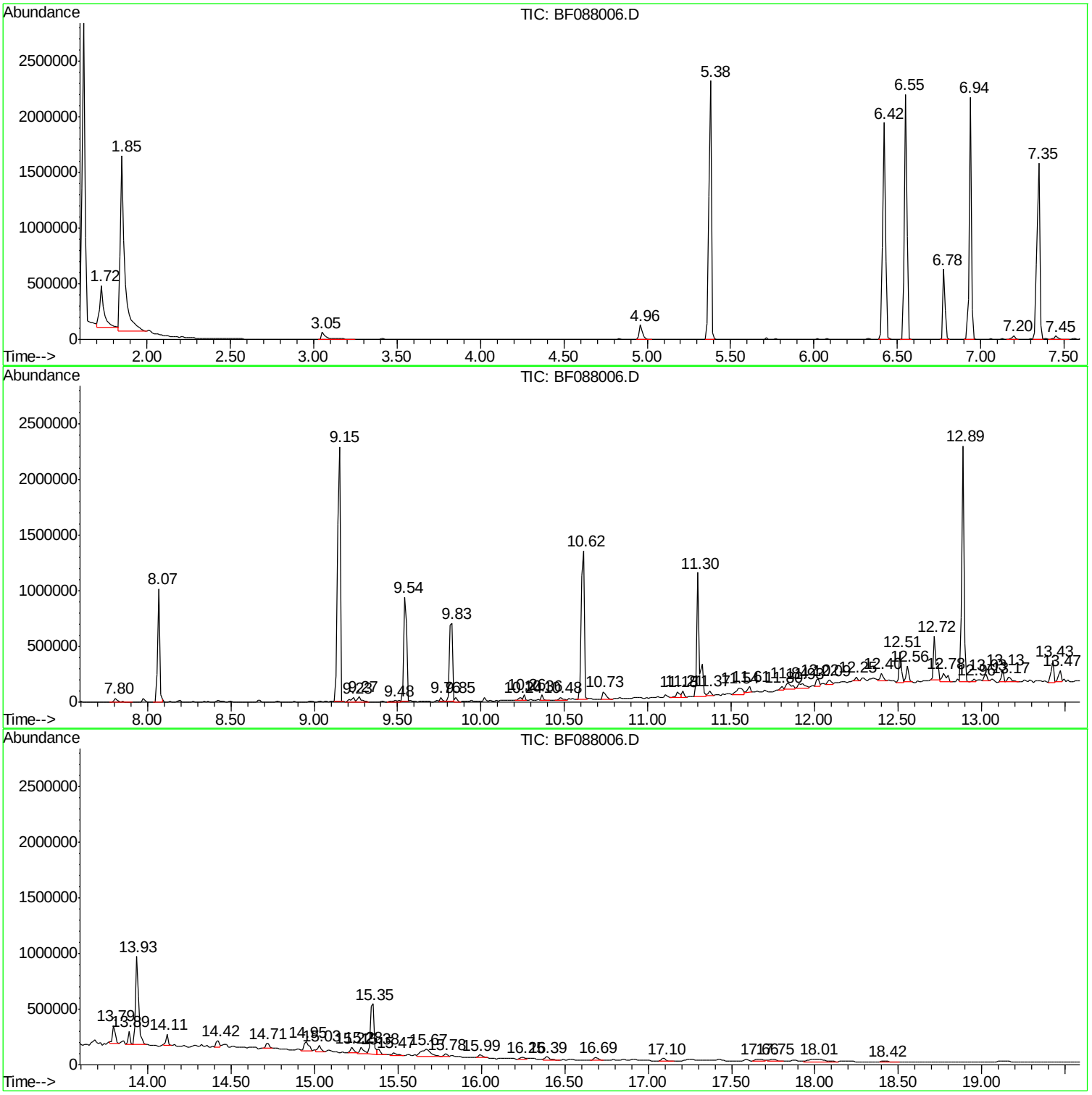
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

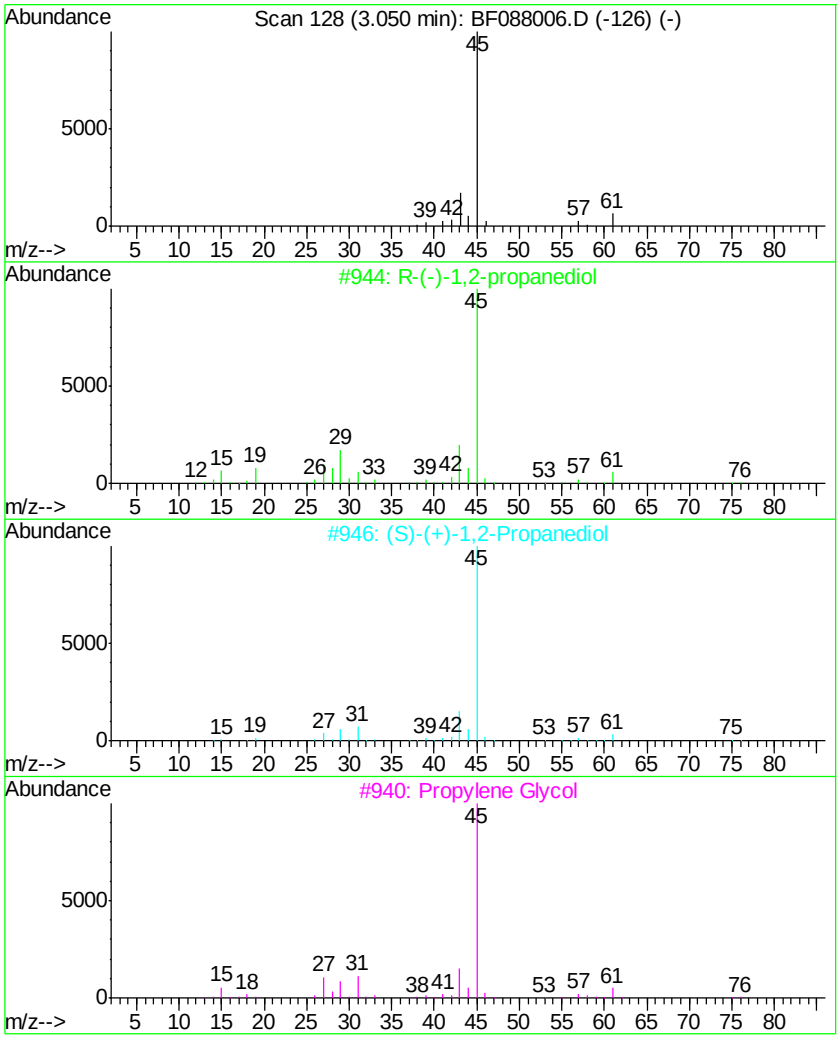
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

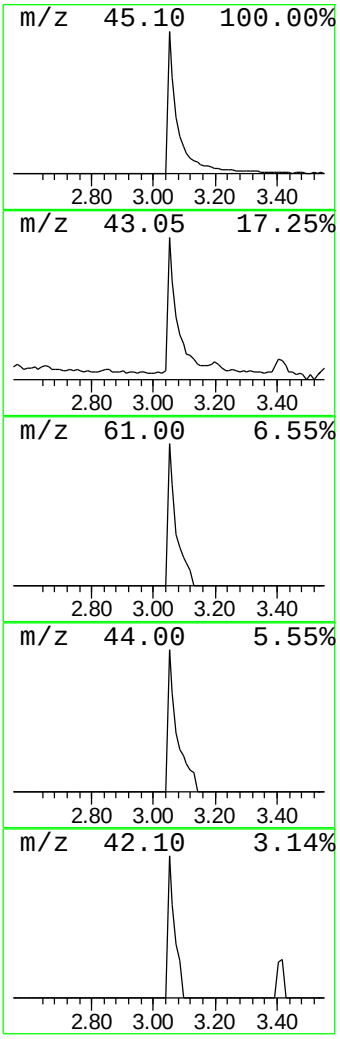
****

Peak Number 3 R-(-)-1,2-propanediol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.05	4.25 ng	133432	1,4-Dichlorobenzene-d4	6.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	64
2	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	9
3	Propylene Glycol	76	C3H8O2	000057-55-6	9
4	Isopropyl Alcohol	60	C3H8O	000067-63-0	9
5	Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	002155-30-8	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

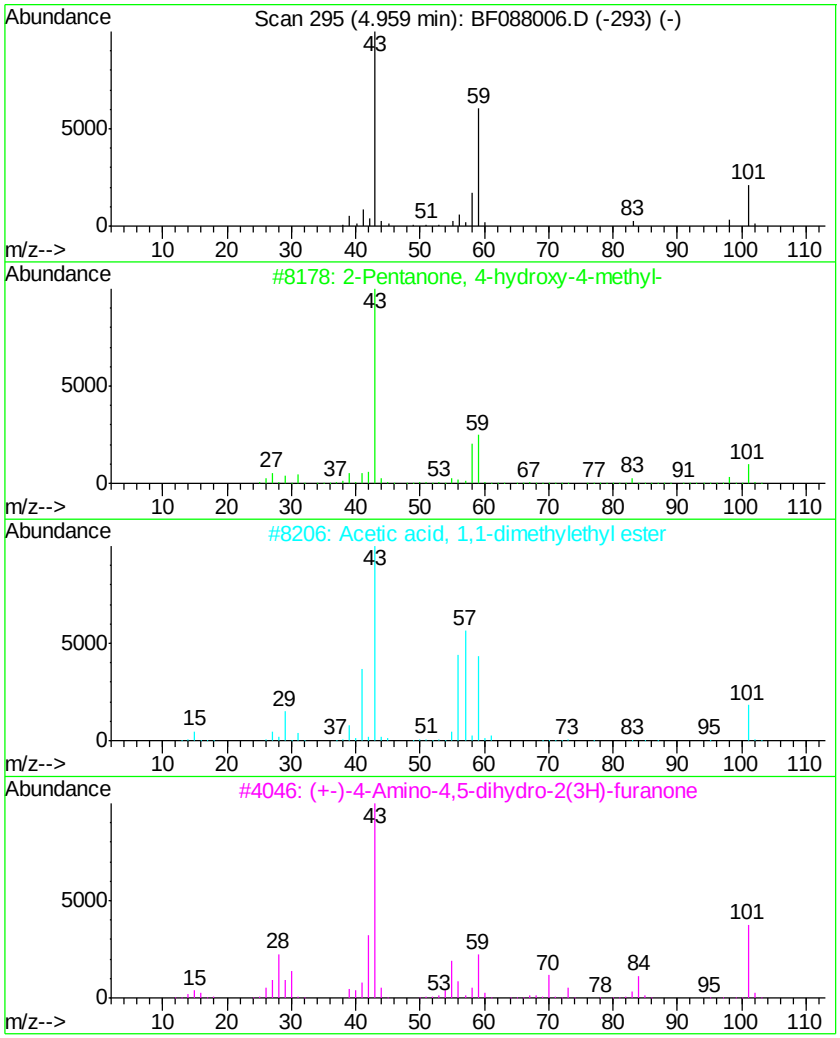
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

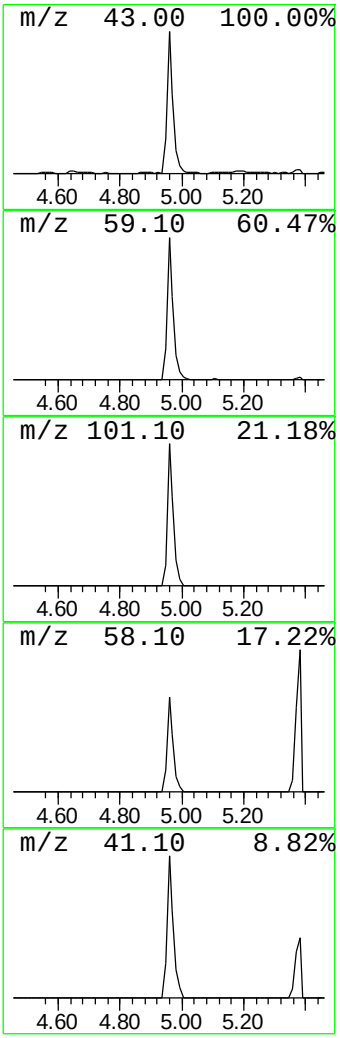
****

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	5.78 ng	181484	1,4-Dichlorobenzene-d4	6.78

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
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	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

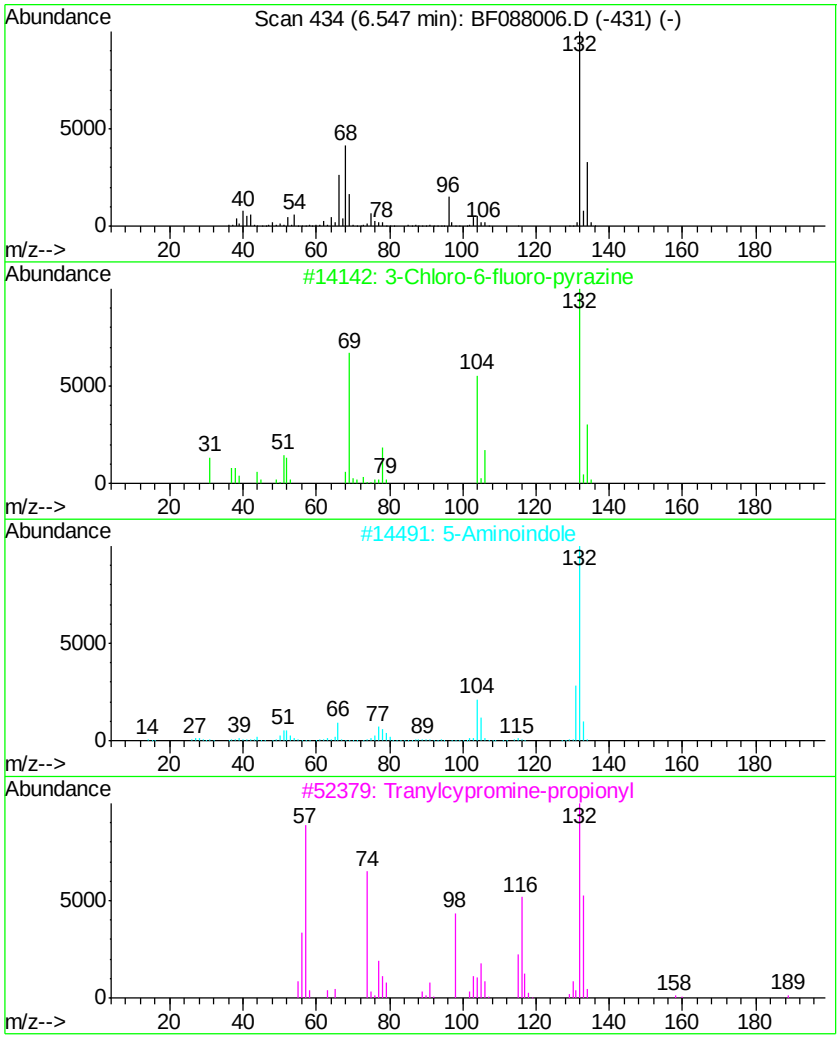
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

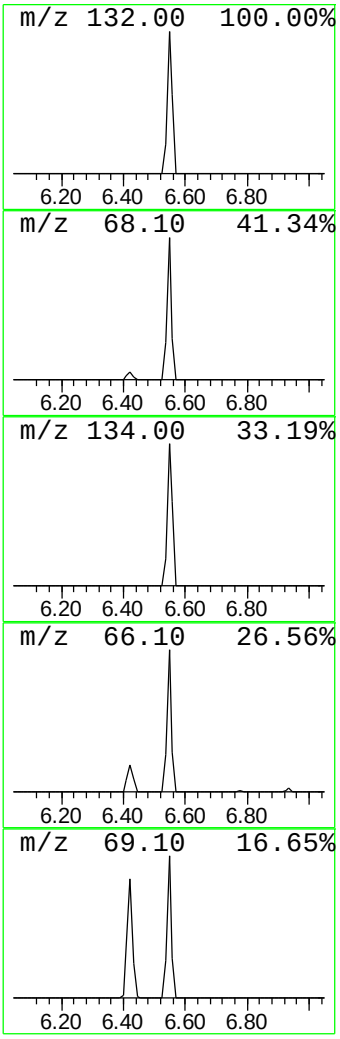
****

Peak Number 5 unknown6.55 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	79.81 ng	2503950	1,4-Dichlorobenzene-d4	6.78

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	12
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9





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Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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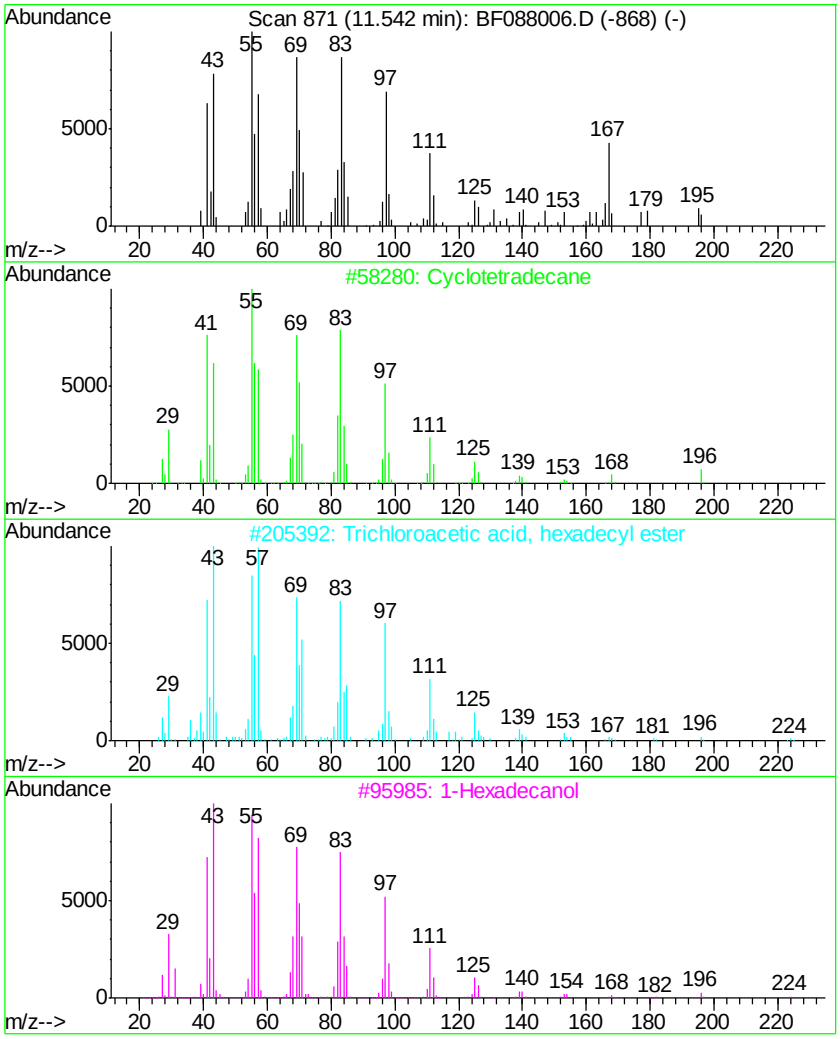
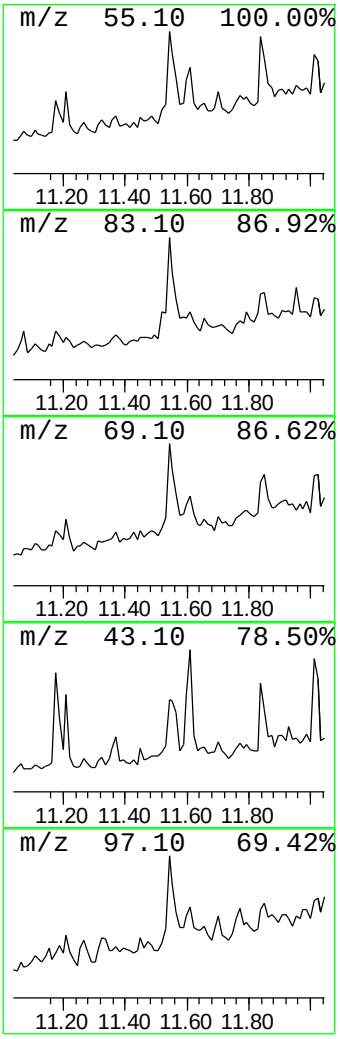
Peak Number 7 Cyclotetradecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.54	2.48 ng	149343	Phenanthrene-d10	11.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	98
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87
3	1-Hexadecanol	242	C16H34O	036653-82-4	87
4	1-Tetradecene	196	C14H28	001120-36-1	86
5	Cyclododecane	168	C12H24	000294-62-2	70





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Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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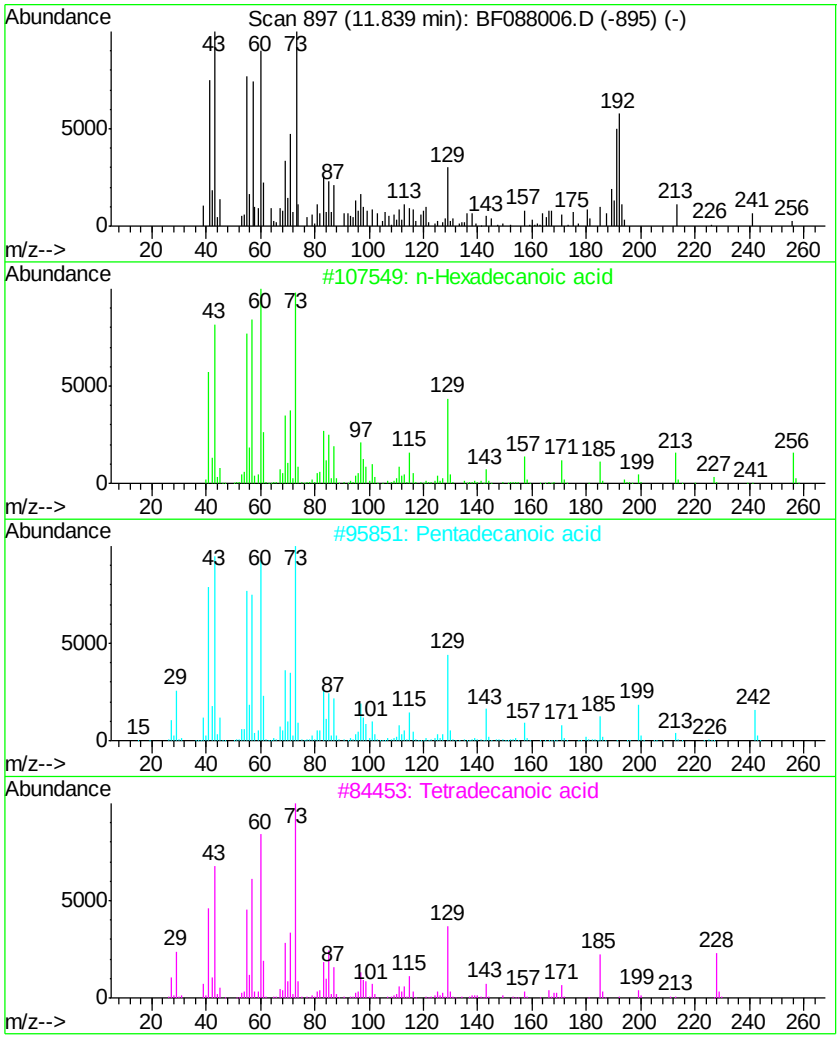
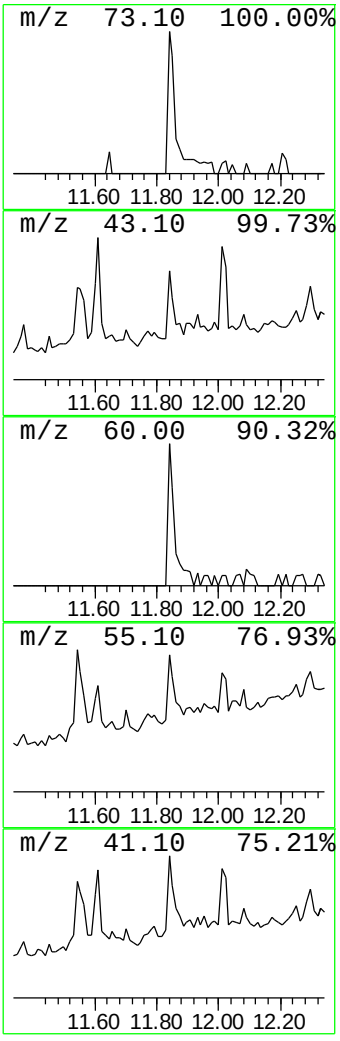
Peak Number 8 n-Hexadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	2.62 ng	157924	Phenanthrene-d10	11.30

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	52
4	n-Decanoic acid	172	C10H20O2	000334-48-5	43
5	Heptanoic acid	130	C7H14O2	000111-14-8	27





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Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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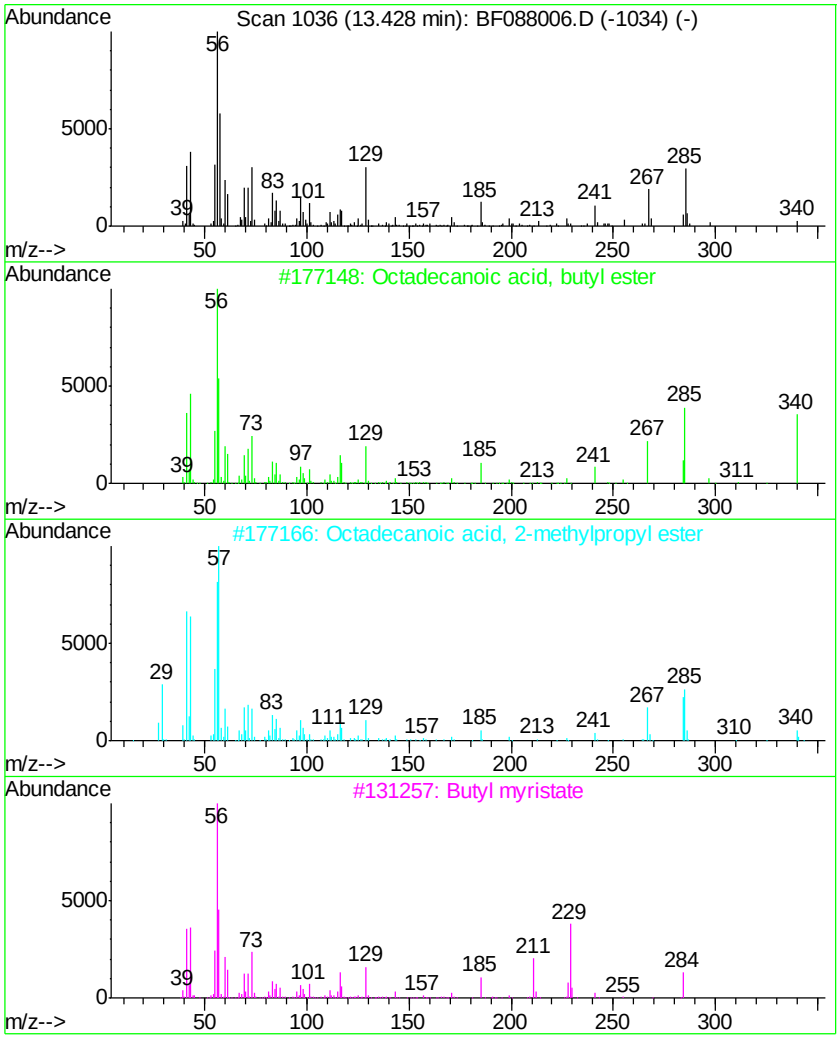
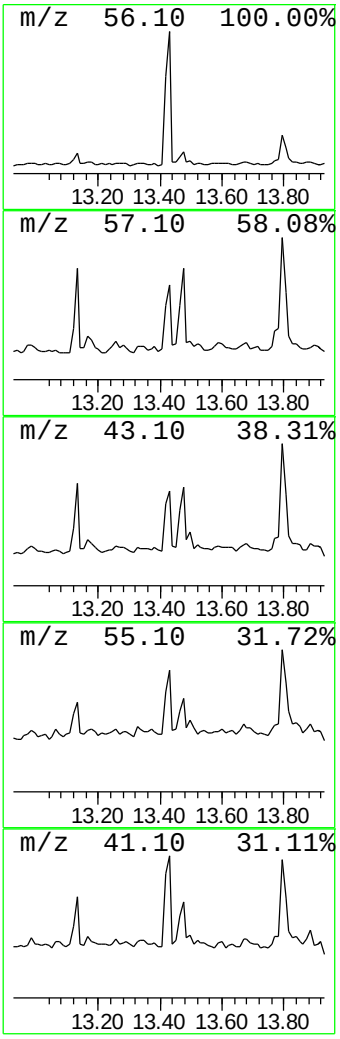
Peak Number 9 Octadecanoic acid, butyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	4.11 ng	215445	Chrysene-d12	13.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	93
2	Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	80
3	Butyl myristate	284	C18H36O2	000110-36-1	50
4	Nipecotic acid	129	C6H11NO2	000498-95-3	47
5	Propan-2-one O-(2-chloro-1-methy...	149	C6H12ClNO	1000186-05-6	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF061416\
Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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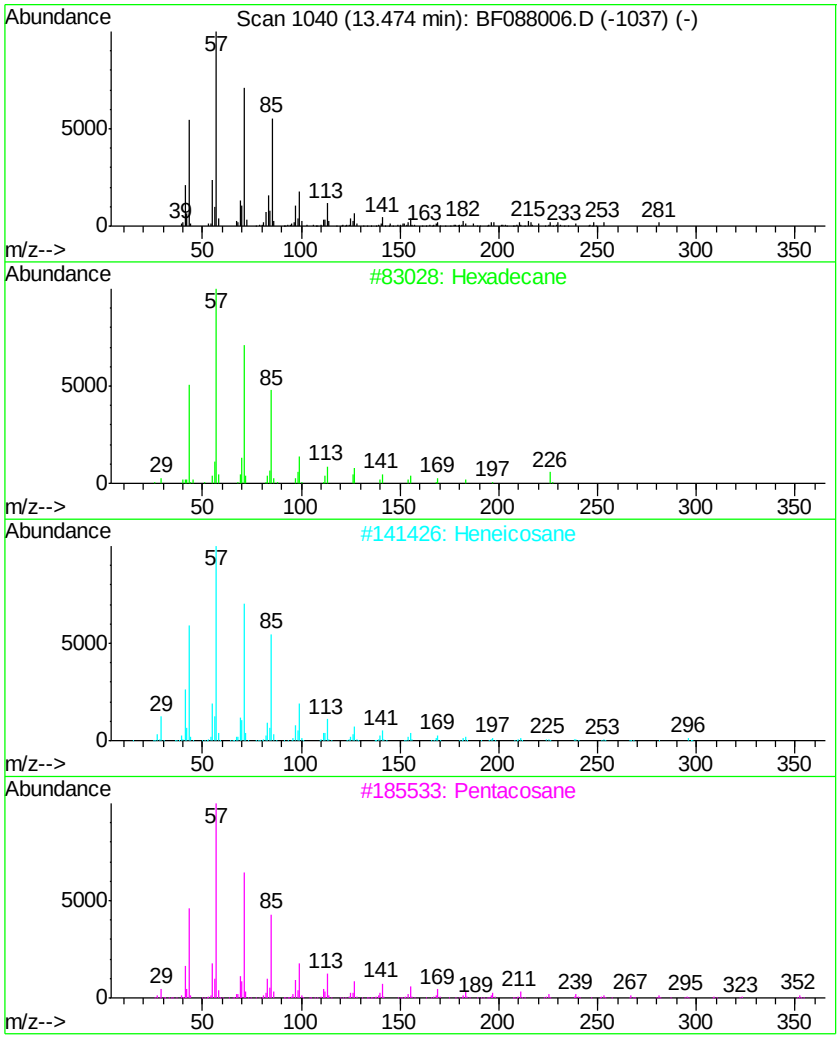
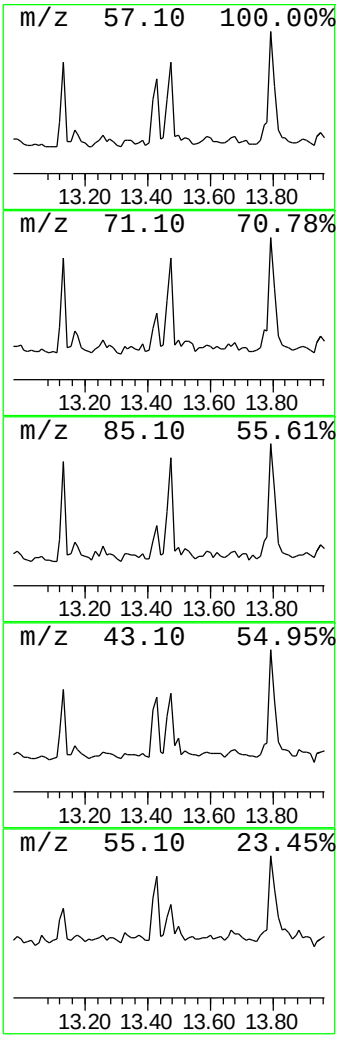
Peak Number 10 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	2.18 ng	114358	Chrysene-d12	13.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Pentacosane	352	C25H52	000629-99-2	90
4	Octacosane	394	C28H58	000630-02-4	87
5	Docosane	310	C22H46	000629-97-0	86





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Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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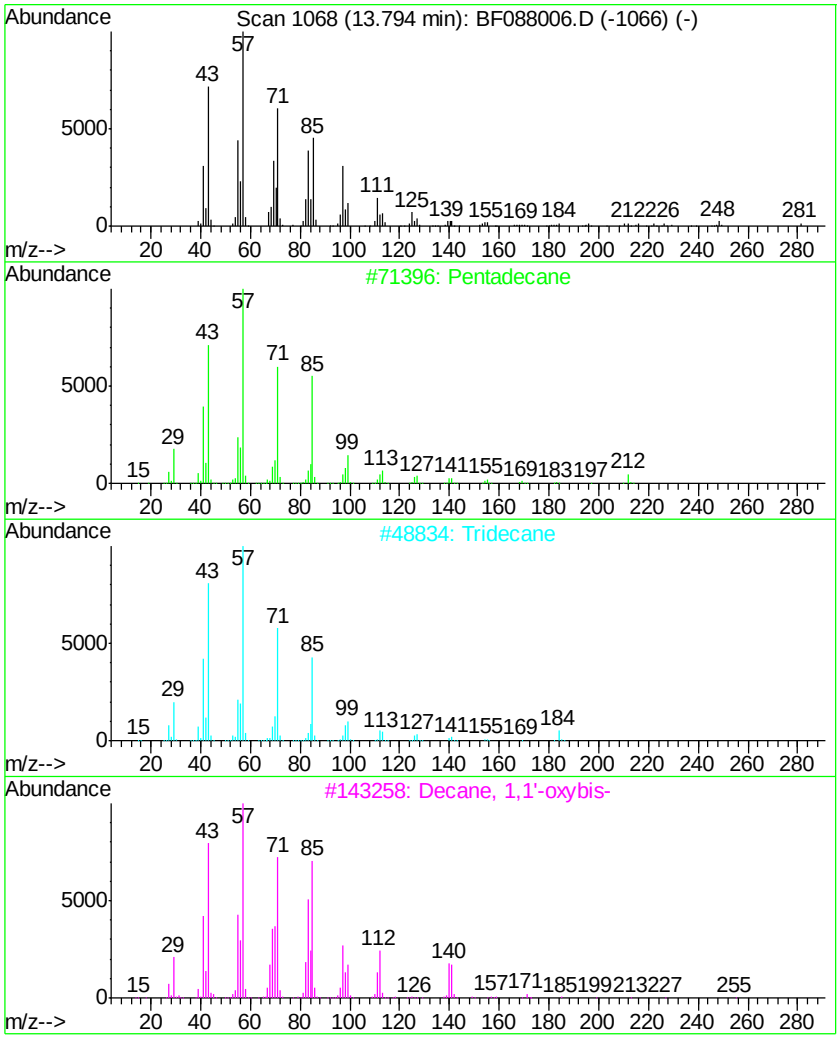
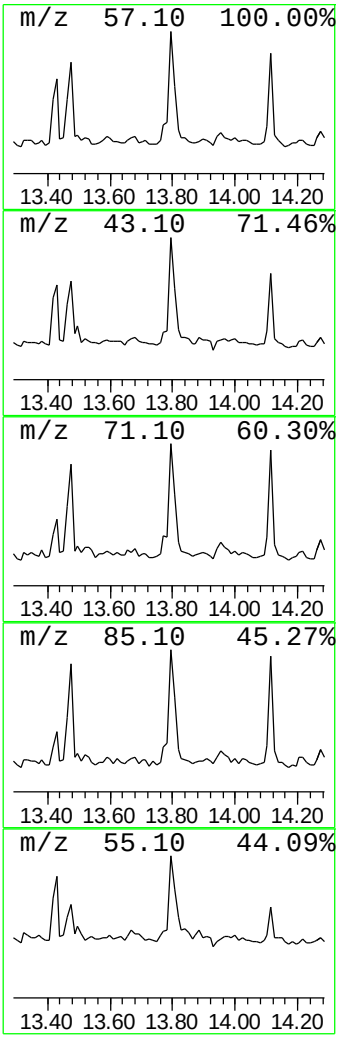
Peak Number 11 Pentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	3.93 ng	206108	Chrysene-d12	13.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	93
2	Tridecane	184	C13H28	000629-50-5	92
3	Decane, 1,1'-oxybis-	298	C20H42O	002456-28-2	87
4	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	87
5	Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	74





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Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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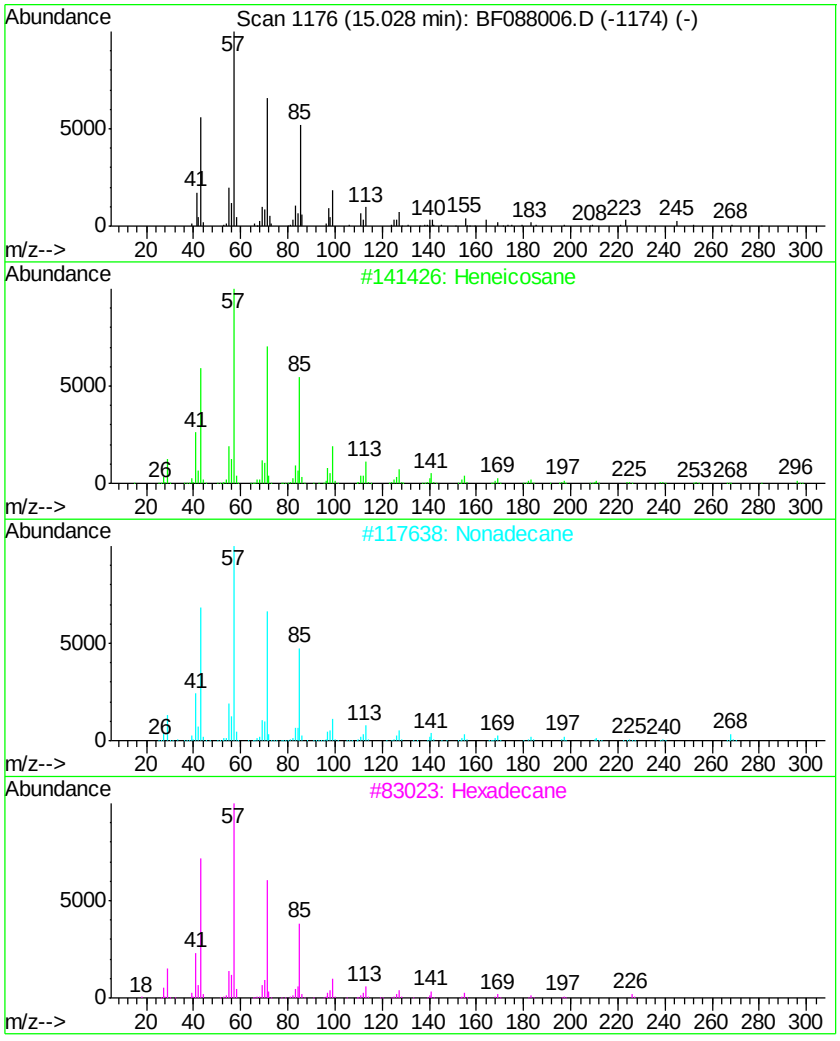
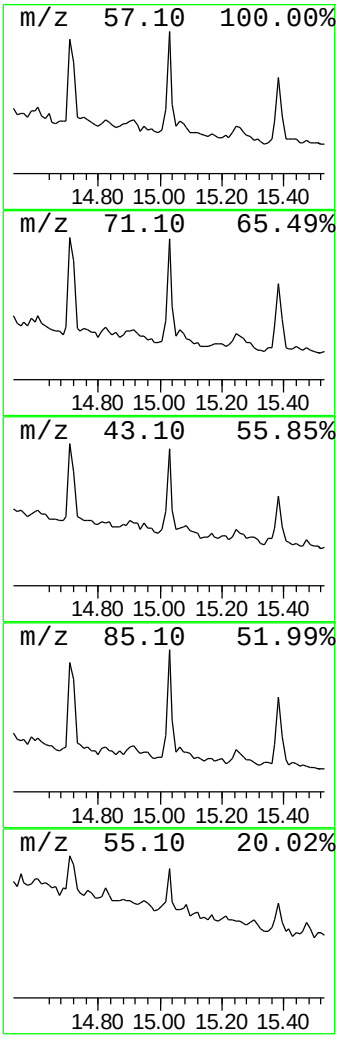
Peak Number 12 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.03	2.23 ng	79324	Perylene-d12	15.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Nonadecane	268	C19H40	000629-92-5	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
5	Octadecane	254	C18H38	000593-45-3	90





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Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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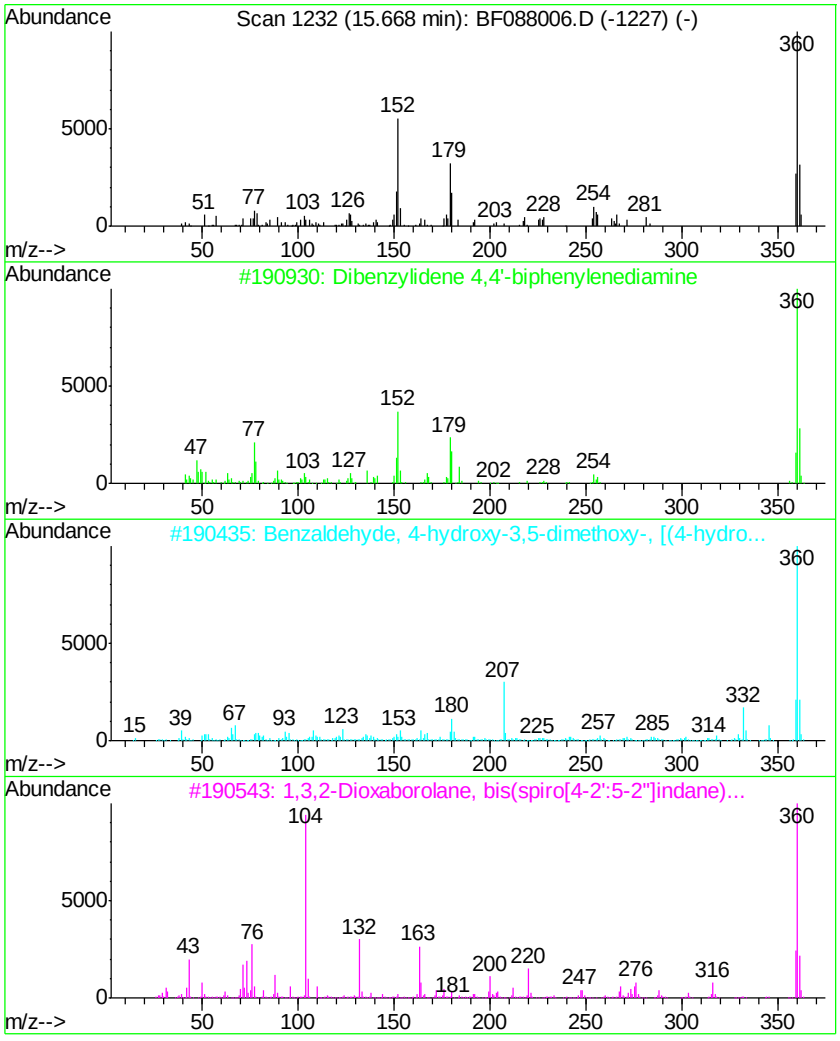
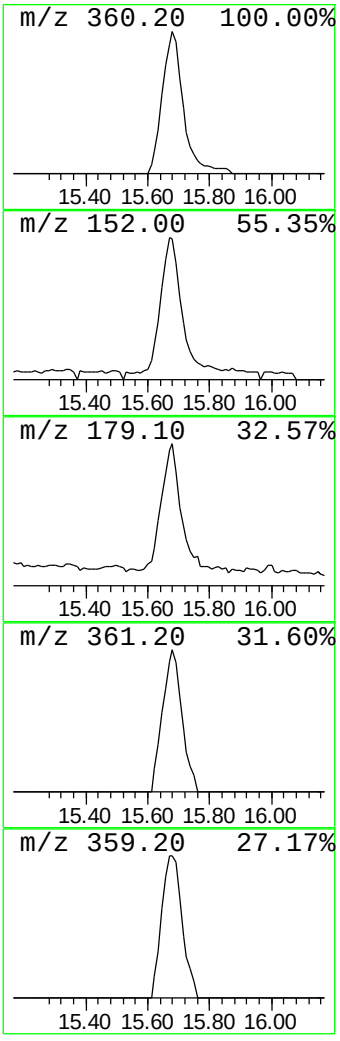
Peak Number 13 Dibenzylidene 4,4'-biphenyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	7.62 ng	270845	Perylene-d12	15.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzylidene 4,4'-biphenylenedi...	360	C26H20N2	006311-48-4	91
2	Benzaldehyde, 4-hydroxy-3,5-dime...	360	C18H20N2O6	014414-32-5	47
3	1,3,2-Dioxaborolane, bis(spiro[4...	360	C20H13BO6	1000150-23-7	47
4	1,10-Phenanthroline, 2,9-dimethy...	360	C26H20N2	004733-39-5	43
5	N,N'-di-2-Naphthyl-p-phenylenedi...	360	C26H20N2	000093-46-9	43





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Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
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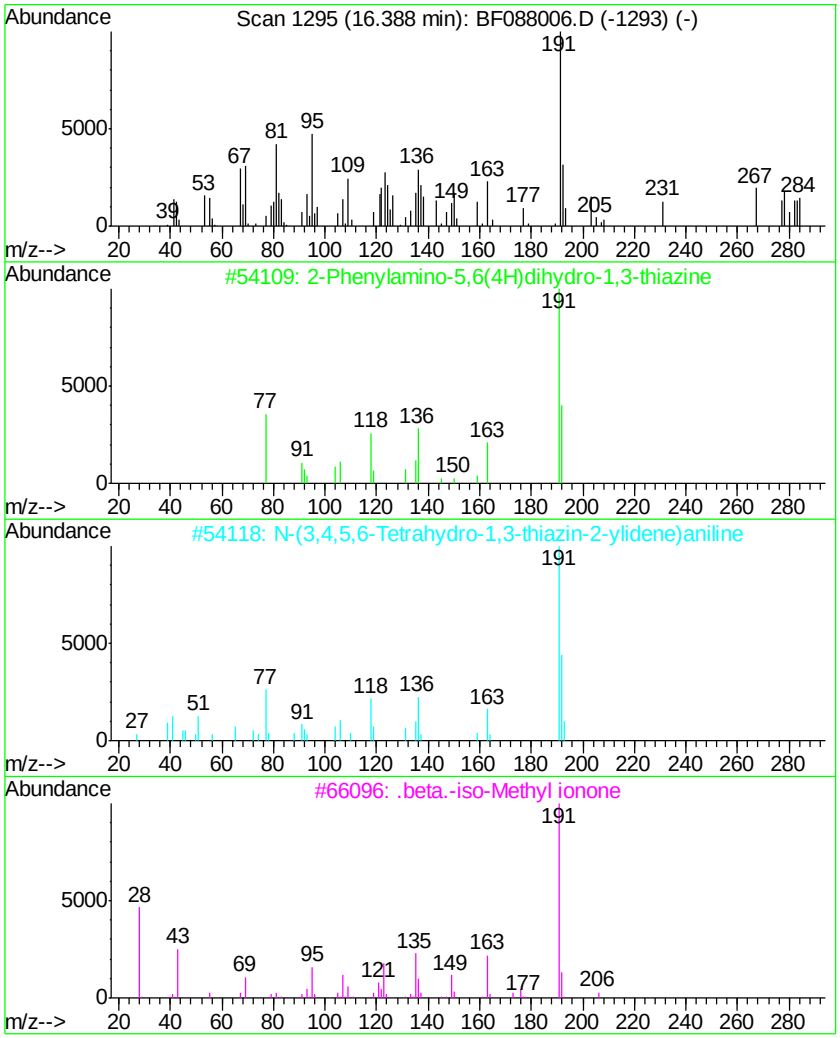
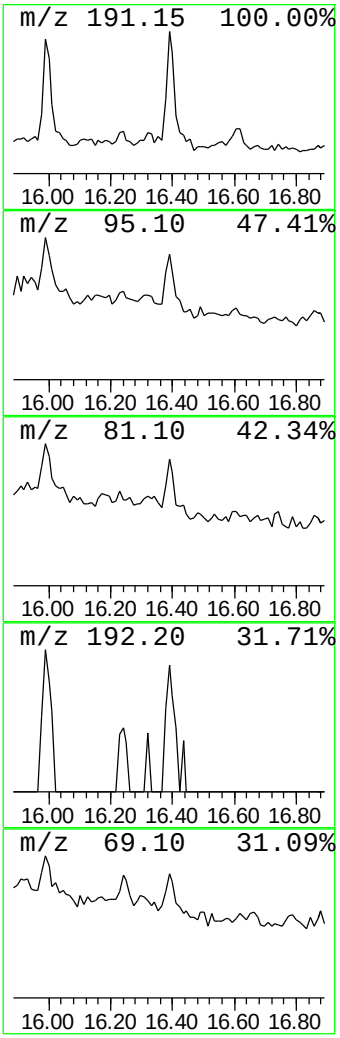
Peak Number 14 unknown16.39 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.39	2.08 ng	73983	Perylene-d12	15.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenylamino-5,6(4H)dihydro-1,3...	192	C10H12N2S	1000147-35-4	38
2	N-(3,4,5,6-Tetrahydro-1,3-thiazi...	192	C10H12N2S	003420-40-4	35
3	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	32
4	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	32
5	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	27





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Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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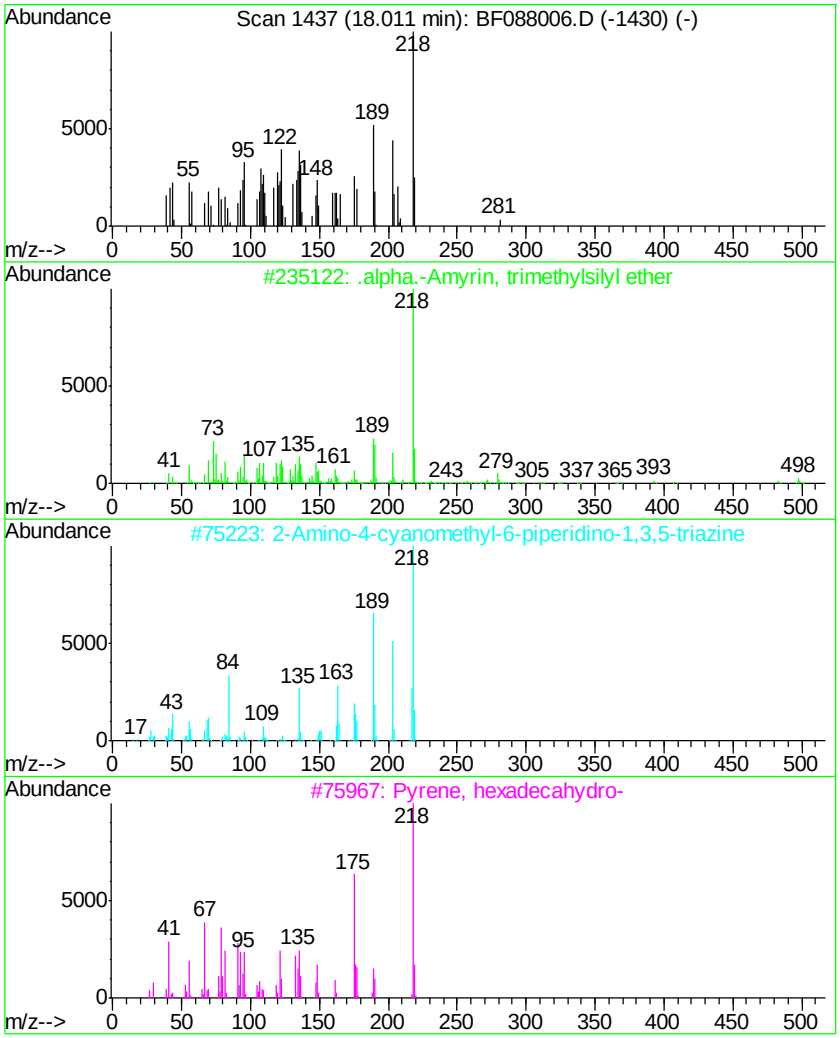
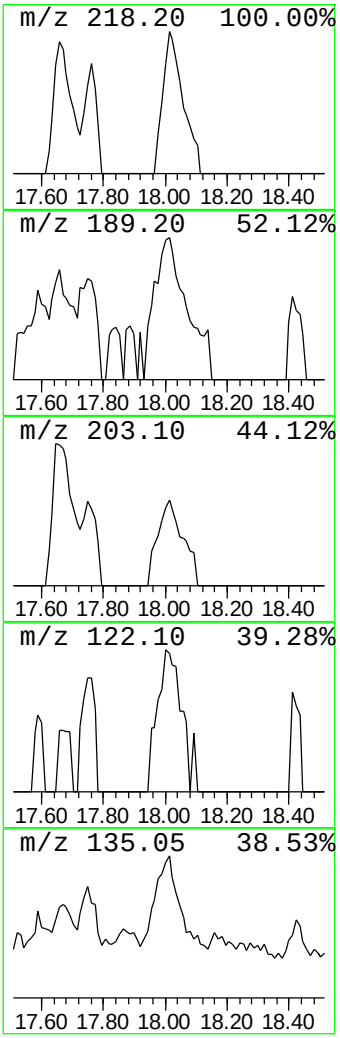
Peak Number 15 unknown18.01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.01	3.78 ng	134425	Perylene-d12	15.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Amyrin, trimethylsilyl e...	498	C33H58OSi	1000374-76-6	43
2	2-Amino-4-cyanomethyl-6-piperidi...	218	C10H14N6	1000241-05-9	43
3	Pyrene, hexadecahydro-	218	C16H26	002435-85-0	43
4	D-Norandrostan-16-one, (5.alpha.)-	260	C18H28O	032319-06-5	38
5	3-Amino-5-piperonyl-1,2,4-triazole	218	C10H10N4O2	014730-92-8	37





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061416\
Data File : BF088006.D
Acq On : 14 Jun 2016 19:01
Operator : UM/SJ
Sample : H3437-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : U6-B1(6-12)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
R-(-)-1,2-propane...	3.05	4.3	ng	133432	1	6.78	627514	20.0
2-Pentanone, 4-hy...	4.96	5.8	ng	181484	1	6.78	627514	20.0
unknown6.55	6.55	79.8	ng	2503950	1	6.78	627514	20.0
Cyclotetradecane	11.54	2.5	ng	149343	4	11.30	1204440	20.0
n-Hexadecanoic acid	11.84	2.6	ng	157924	4	11.30	1204440	20.0
Octadecanoic acid...	13.43	4.1	ng	215445	5	13.93	1049330	20.0
Hexadecane	13.47	2.2	ng	114358	5	13.93	1049330	20.0
Pentadecane	13.79	3.9	ng	206108	5	13.93	1049330	20.0
Heneicosane	15.03	2.2	ng	79324	6	15.35	711161	20.0
Dibenzylidene 4,4...	15.67	7.6	ng	270845	6	15.35	711161	20.0
unknown16.39	16.39	2.1	ng	73983	6	15.35	711161	20.0
unknown18.01	18.01	3.8	ng	134425	6	15.35	711161	20.0