LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

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-BF063016.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	1.667	6	7	12	rVV	83237	120368	6.75%	0.737%
2	1.793	16	18	25	rVV	491593	667015	37.43%	4.086%
3	1.896	25	27	40	rVB	45676	104968	5.89%	0.643%
4	2.867	111	112	121	rBV	19781	42889	2.41%	0.263%
5	4.856	284	286	291	rBV	123970	195731	10.98%	1.199%
6	5.302	323	325	328	rBV	1245160	1470993	82.54%	9.011%
7	6.353	414	417	420	rVB	1181420	1502507	84.30%	9.204%
8	6.479	425	428	430	rVB	1621630	1777462	99.73%	10.888%
9	6.707	446	448	449	rBV	346991	387943	21.77%	2.376%
10	6.856	459	461	464	rVB	1037511	1225934	68.79%	7.510%
11	7.050	475	478	481	rBV2	11356	23604	1.32%	0.145%
12	7.268	494	497	499	rBV	843519	941952	52.85%	5.770%
13	7.988	558	560	563	rBV	574174	508488	28.53%	3.115%
14	8.616	612	615	617	rBV	145914	143222	8.04%	0.877%
15	9.073	652	655	657	rBV	1699543	1782236	100.00%	10.917%
16	9.405	680	684	687	rBV	11823	21295	1.19%	0.130%
17	9.462	687	689	691	rVV	266715	220951	12.40%	1.353%
18	9.553	693	697	699	rVV	14386	25569	1.43%	0.157%
19	9.599	699	701	703	rVB	26296	19717	1.11%	0.121%
20	9.736	711	713	715	rBV	497043	457936	25.69%	2.805%
21	9.862	723	724	727	rVB2	50464	50981	2.86%	0.312%
22	10.239	754	757	759	rBV	20551	23664	1.33%	0.145%
23	10.525	780	782	784	rBV	839486	768917	43.14%	4.710%
24	10.651	792	793	797	rVB2	24433	46607	2.62%	0.285%
25	10.971	815	821	824	rBV	85607	133062	7.47%	0.815%
26	11.096	831	832	834	rBV	32758	42167	2.37%	0.258%
27	11.211	840	842	846	rVB	368997	374084	20.99%	2.291%
28	11.291	846	849	851	rBV	26692	29857	1.68%	0.183%
29	11.451	861	863	868	rVB2	14345	26617	1.49%	0.163%
30	11.531	868	870	875	rVB	20162	27392	1.54%	0.168%
31	11.759	887	890	891	rBV2	35519	39602	2.22%	0.243%
32	11.817	891	895	901	rVB2	45559	84097	4.72%	0.515%
33	11.931	901	905	909	rBV2	11030	18175	1.02%	0.111%
34	12.262	931	934	936	rBV	23763	26440	1.48%	0.162%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

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-BF063016.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.422	945	948	950	rBV	54099	57627	3.23%	0.353%
36	12.639	965	967	969	rBV	54850	65051	3.65%	0.398%
37	12.799	979	981	983	rVB	1244139	1086575	60.97%	6.656%
38	12.948	990	994	998	rVB2	12247	29289	1.64%	0.179%
39	13.040	998	1002	1004	rBV2	15602	34042	1.91%	0.209%
40	13.360	1027	1030	1031	rBV3	10666	22717	1.27%	0.139%
41	13.474	1038	1040	1045	rBV	17328	22391	1.26%	0.137%
42	13.588	1048	1050	1051	rBV2	17360	23247	1.30%	0.142%
43	13.702	1057	1060	1064	rBV2	114756	123884	6.95%	0.759%
44	13.840	1069	1072	1075	rBV	464459	527063	29.57%	3.229%
45	14.262	1107	1109	1110	rBV2	16696	19619	1.10%	0.120%
46	14.331	1113	1115	1121	rVV4	26325	70483	3.95%	0.432%
47	14.605	1137	1139	1140	rBV	16804	30232	1.70%	0.185%
48	14.754	1150	1152	1153	rBV2	17285	21840	1.23%	0.134%
49	14.834	1156	1159	1162	rVV	102149	173724	9.75%	1.064%
50	14.925	1166	1167	1171	rVV	33973	44380	2.49%	0.272%
51	15.097	1180	1182	1185	rVB	54446	68166	3.82%	0.418%
52	15.154	1185	1187	1190	rBV	50936	59509	3.34%	0.365%
53	15.211	1190	1192	1195	rBV	286513	339182	19.03%	2.078%
54	15.645	1228	1230	1232	rBV	22178	30161	1.69%	0.185%
55	16.491	1302	1304	1307	rBV2	23783	45747	2.57%	0.280%
56	16.651	1316	1318	1321	rBV3	25619	48723	2.73%	0.298%
57	16.880	1336	1338	1342	rVB	17242	29981	1.68%	0.184%
58	17.703	1408	1410	1413	rVB4	18736	18782	1.05%	0.115%

Sum of corrected areas: 16324857

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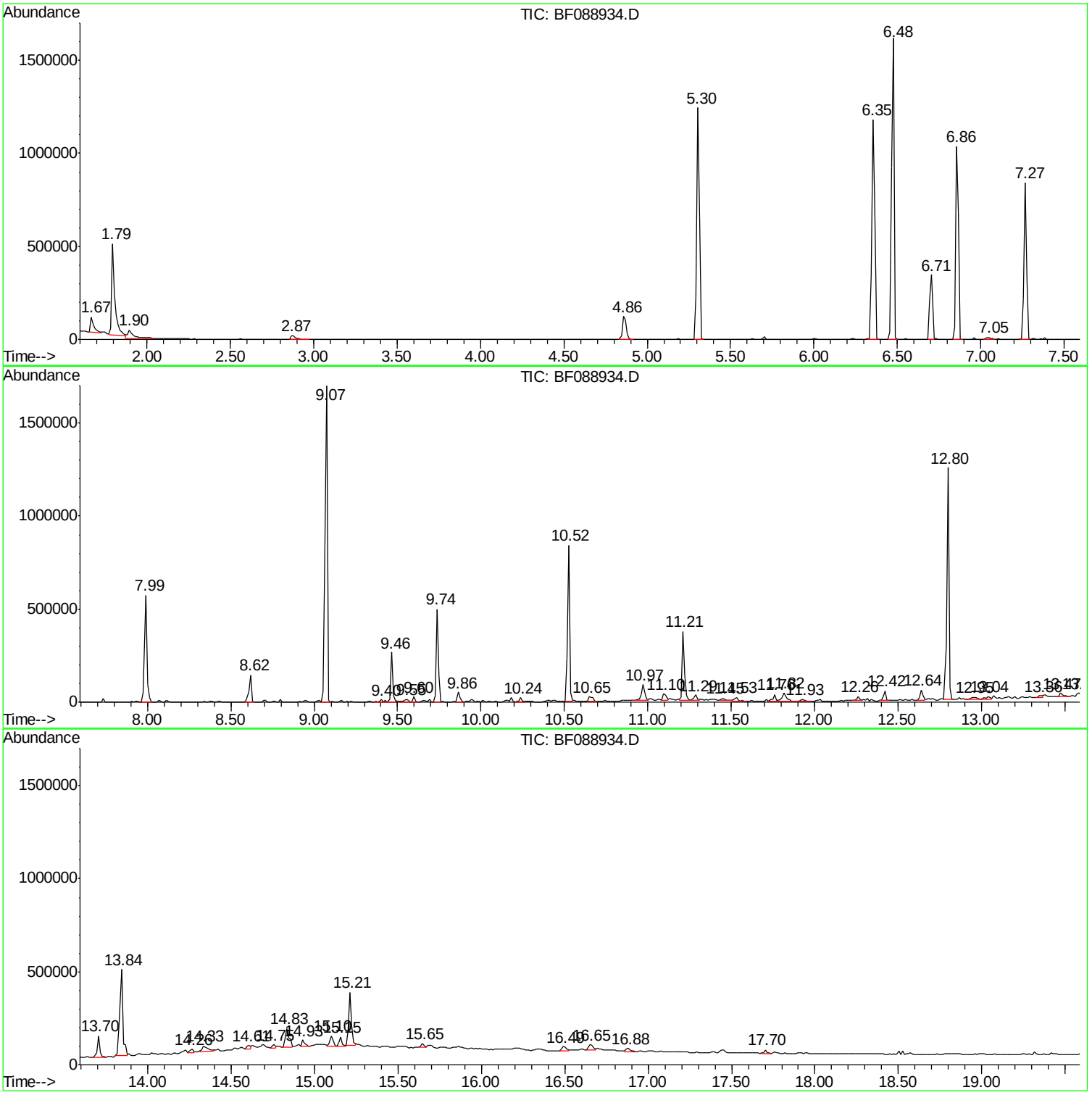
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
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Operator : UM/SJ
Sample : H3934-15
Misc :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

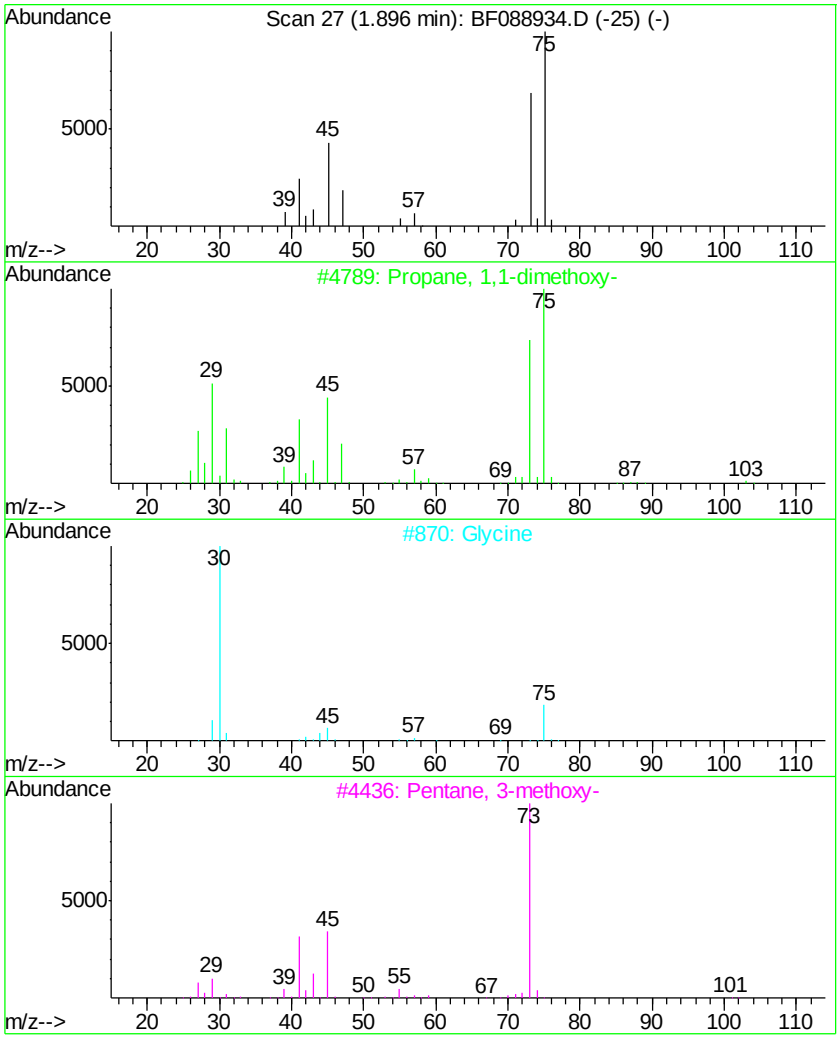
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

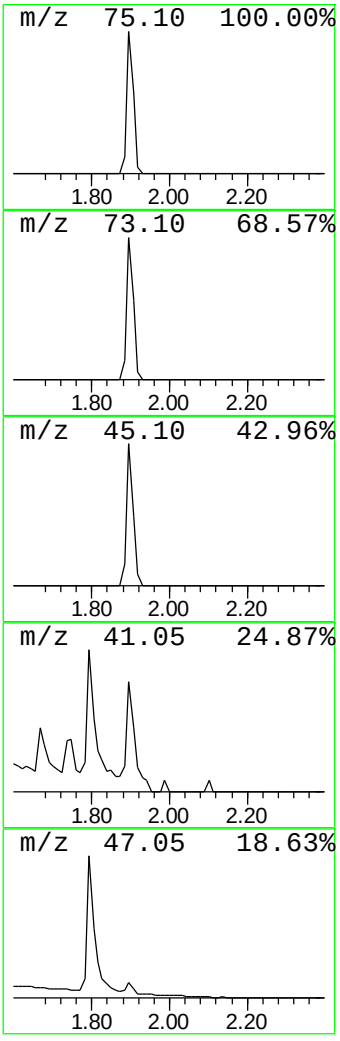
****

Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.90	5.41 ng	104968	1,4-Dichlorobenzene-d4	6.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2	Glycine	75	C2H5NO2	000056-40-6	9
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
4	1-Propene-1-thiol	74	C3H6S	000925-89-3	5
5	Methane, nitroso-	45	CH3NO	000865-40-7	4





Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

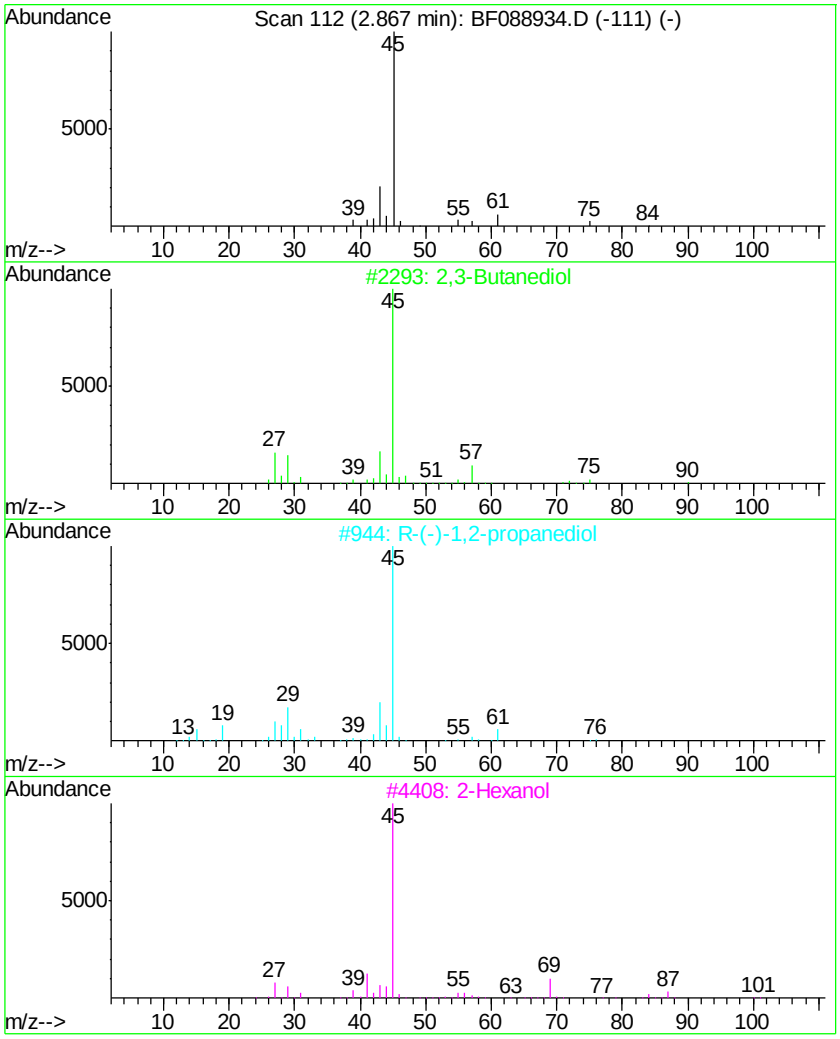
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

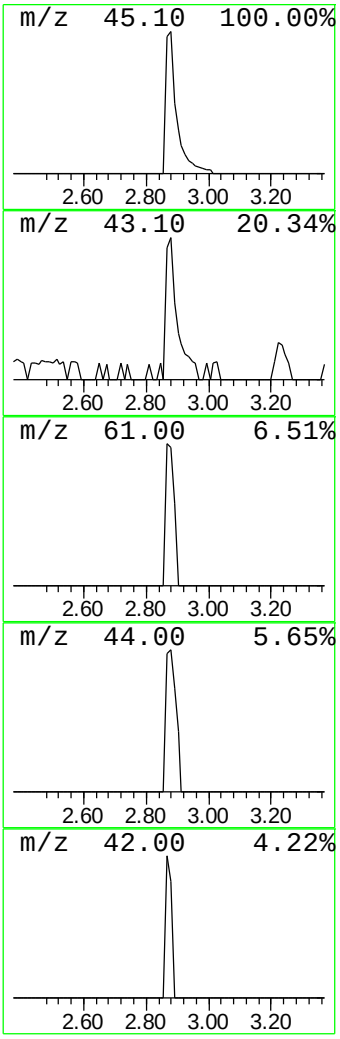
****

Peak Number 4 unknown2.87 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	2.21 ng	42889	1,4-Dichlorobenzene-d4	6.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Butanediol	90	C4H10O2	000513-85-9	43
2	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
3	2-Hexanol	102	C6H14O	000626-93-7	40
4	2-Methoxyisopropyl cyanoacetate	157	C7H11NO3	032804-79-8	9
5	Propylene Glycol	76	C3H8O2	000057-55-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

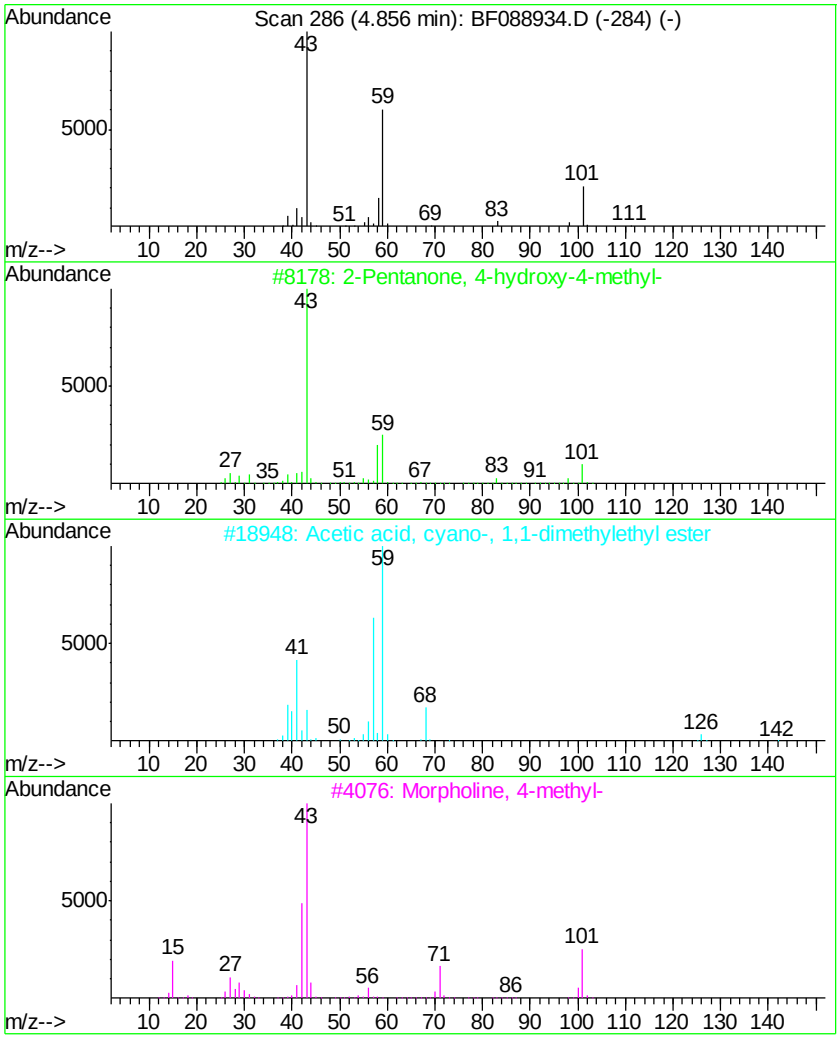
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

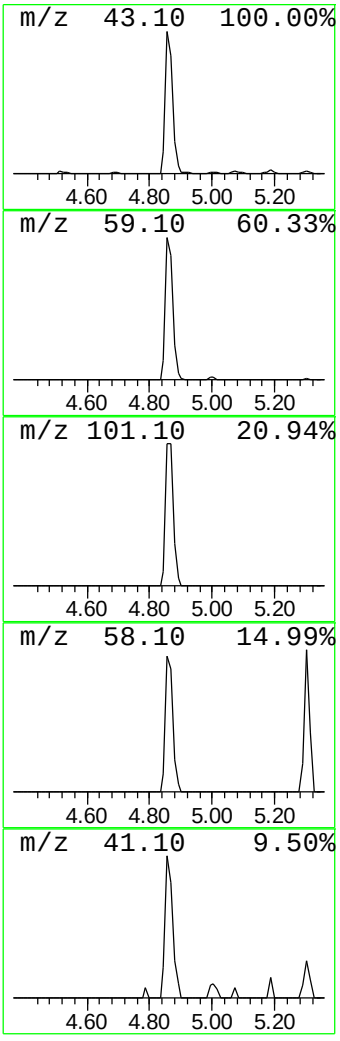
****

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	10.09 ng	195731	1,4-Dichlorobenzene-d4	6.71

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	25
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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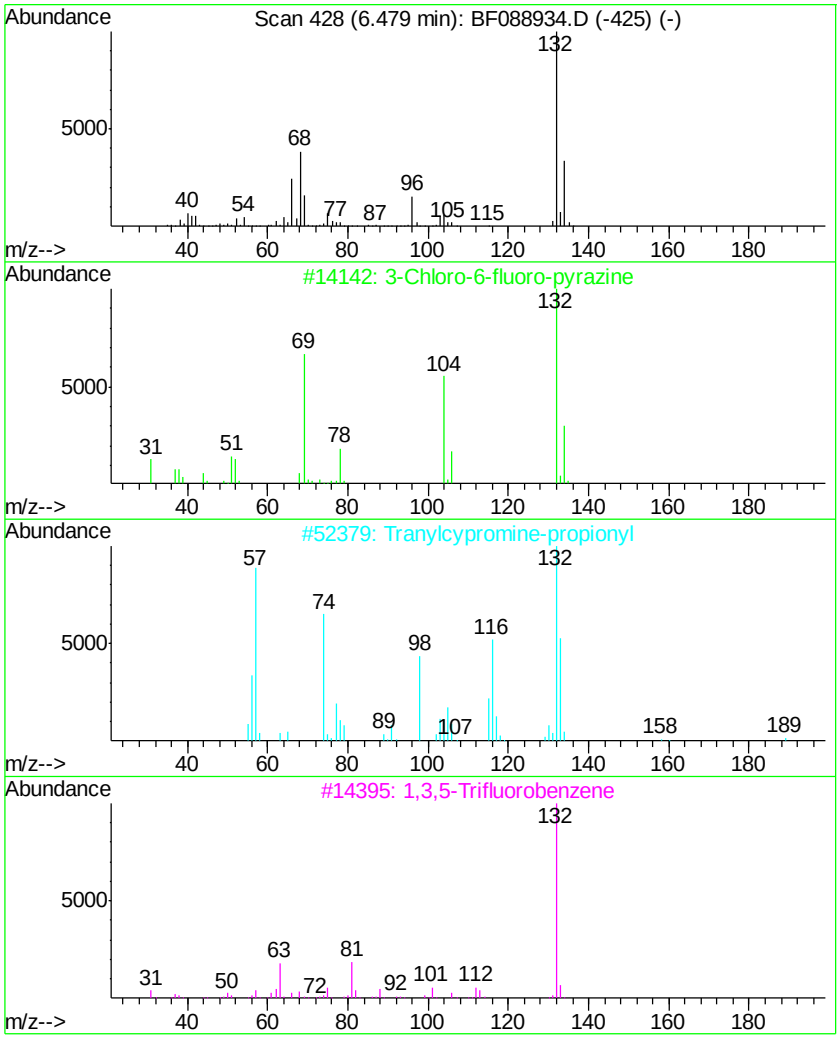
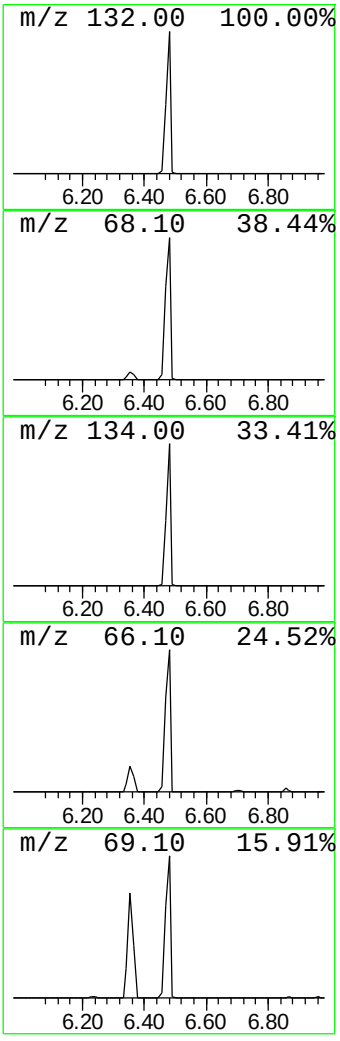
Peak Number 6 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	91.64 ng	1777460	1,4-Dichlorobenzene-d4	6.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9





Library Search Compound Report

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Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
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Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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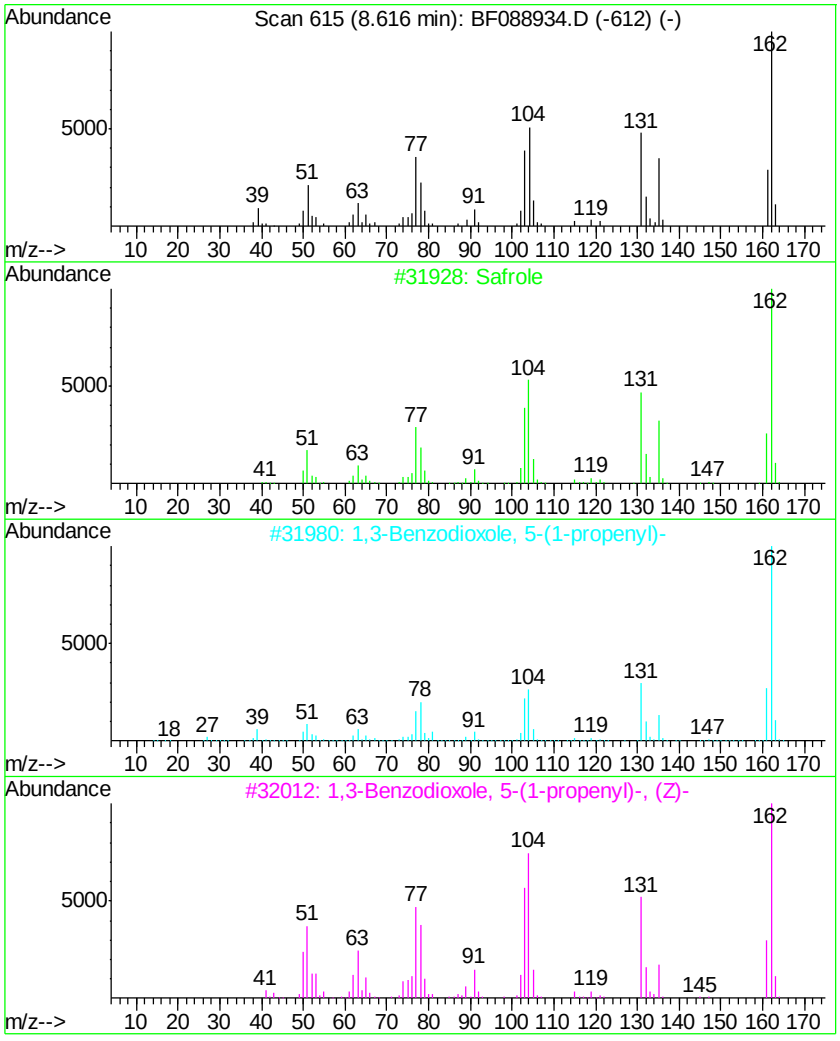
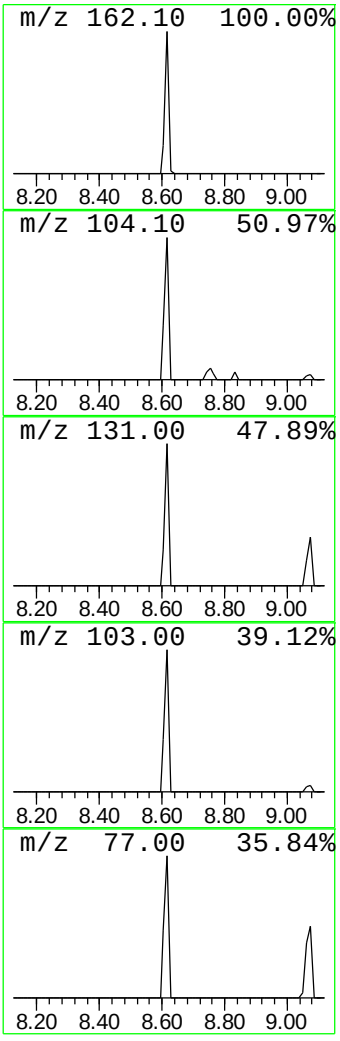
Peak Number 8 Safrole Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.62	5.63 ng	143222	Naphthalene-d8	7.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Safrole	162	C10H10O2	000094-59-7	98
2	1,3-Benzodioxole, 5-(1-propenyl)-	162	C10H10O2	000120-58-1	97
3	1,3-Benzodioxole, 5-(1-propenyl)...	162	C10H10O2	017627-76-8	97
4	Benzene, 1,2-(methylenedioxy)-4-...	162	C10H10O2	004043-71-4	96
5	1H-Isoindole-1,3(2H)-dione, 2-am...	162	C8H6N2O2	001875-48-5	43





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Sample : H3934-15
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ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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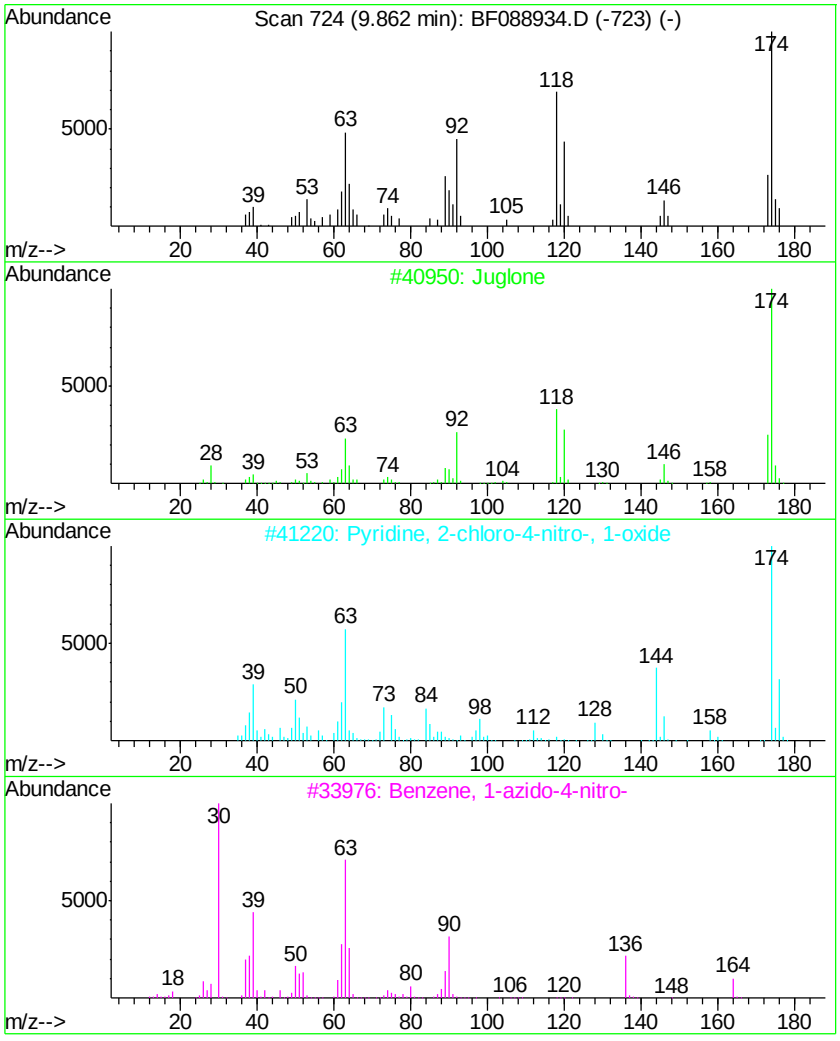
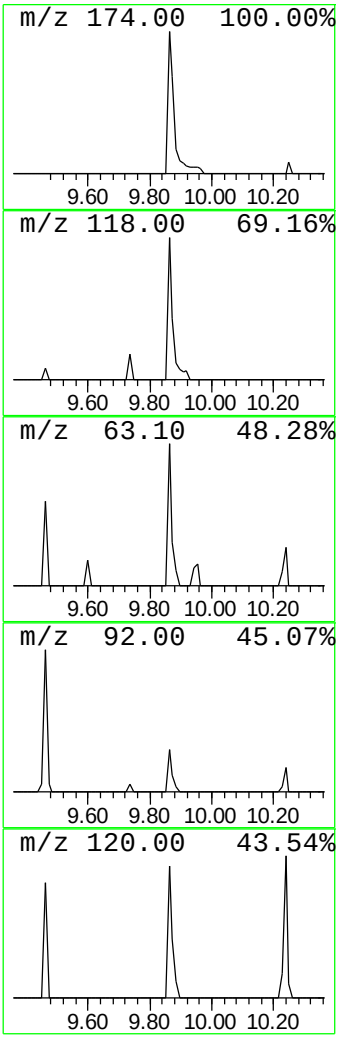
Peak Number 9 Juglone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	2.23 ng	50981	Acenaphthene-d10	9.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Juglone	174	C10H6O3	000481-39-0	97
2	Pyridine, 2-chloro-4-nitro-, 1-o...	174	C5H3ClN2O3	014432-16-7	47
3	Benzene, 1-azido-4-nitro-	164	C6H4N4O2	001516-60-5	25
4	8-Quinolinol, 5-nitroso-	174	C9H6N2O2	003565-26-2	25
5	Isophosphinoline, 1,3-dimethyl-	174	C11H11P	057328-61-7	25





Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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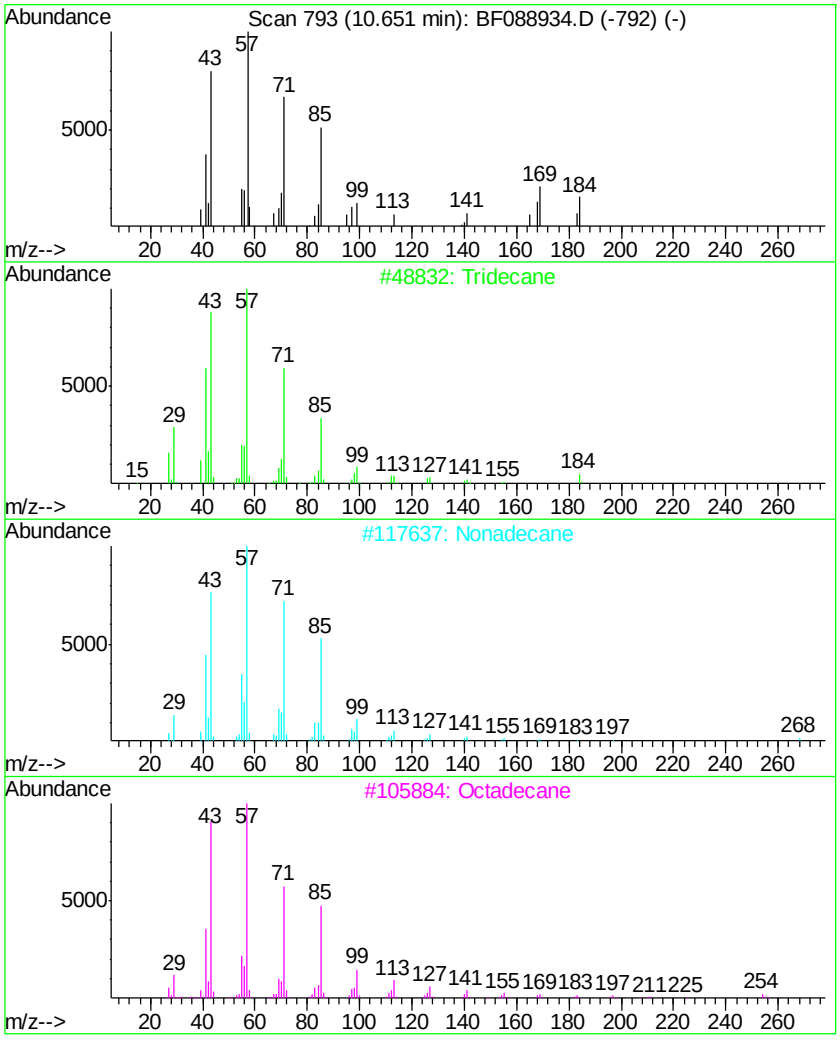
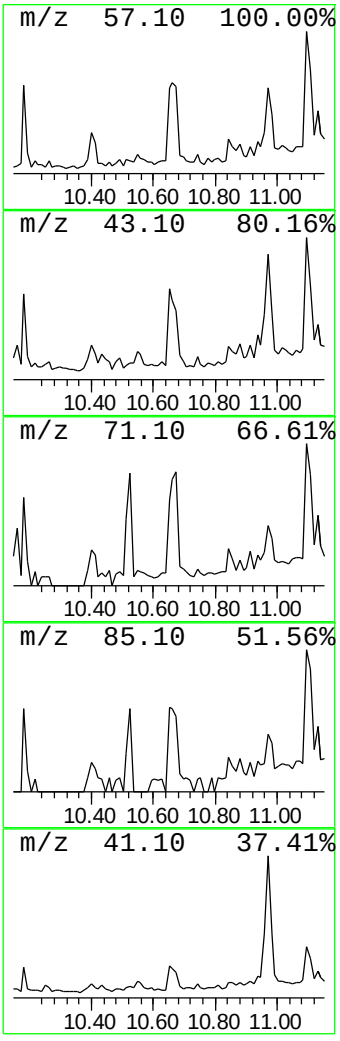
Peak Number 10 Tridecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	2.49 ng	46607	Phenanthrene-d10	11.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	93
2	Nonadecane	268	C19H40	000629-92-5	59
3	Octadecane	254	C18H38	000593-45-3	59
4	Heneicosane	296	C21H44	000629-94-7	53
5	Tetradecane	198	C14H30	000629-59-4	53





Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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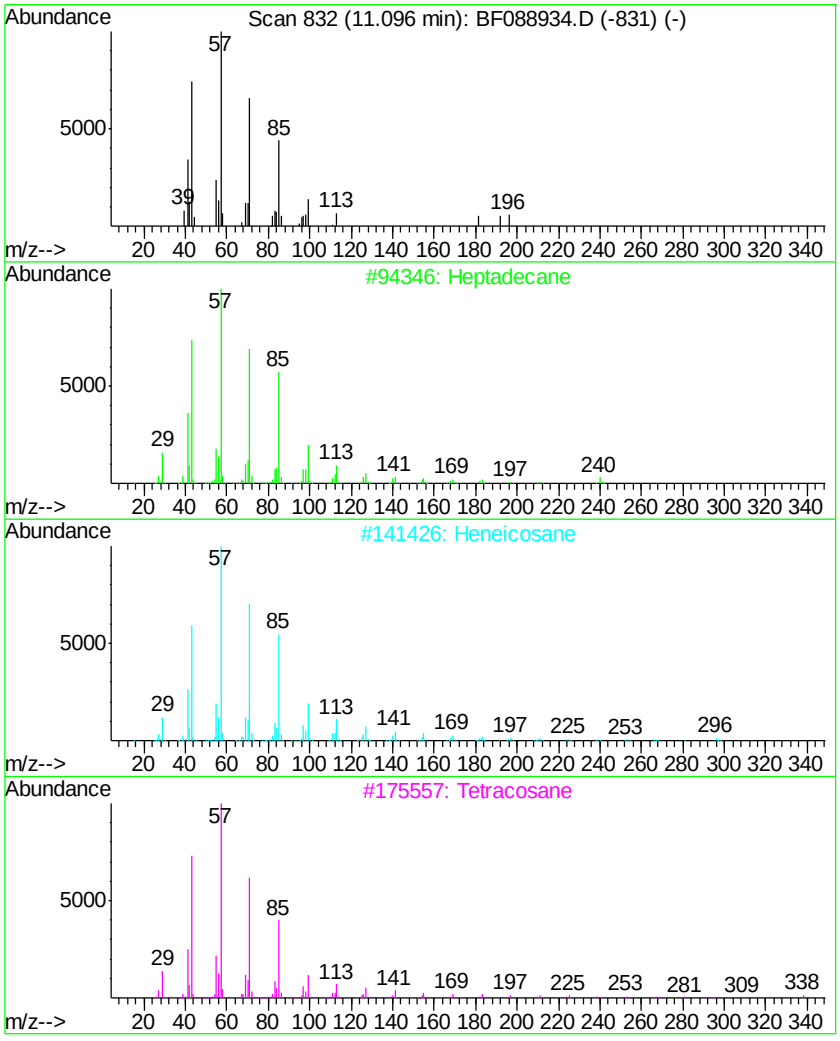
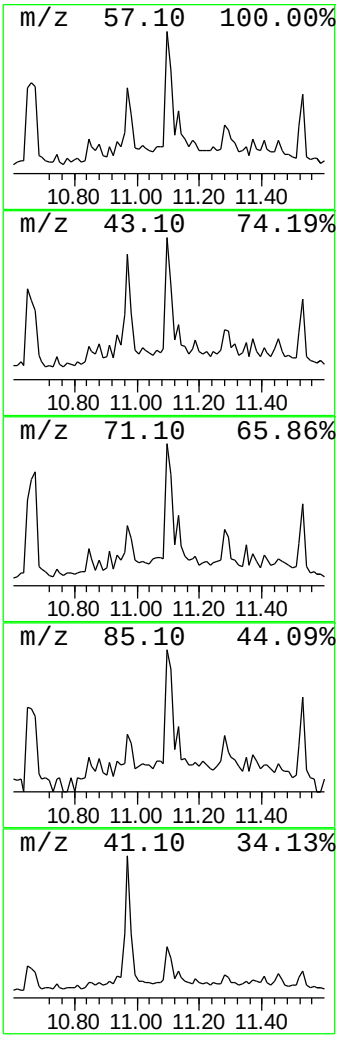
Peak Number 12 Heptadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.25 ng	42167	Phenanthrene-d10	11.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	86
2	Heneicosane	296	C21H44	000629-94-7	86
3	Tetracosane	338	C24H50	000646-31-1	86
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	80
5	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	80





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Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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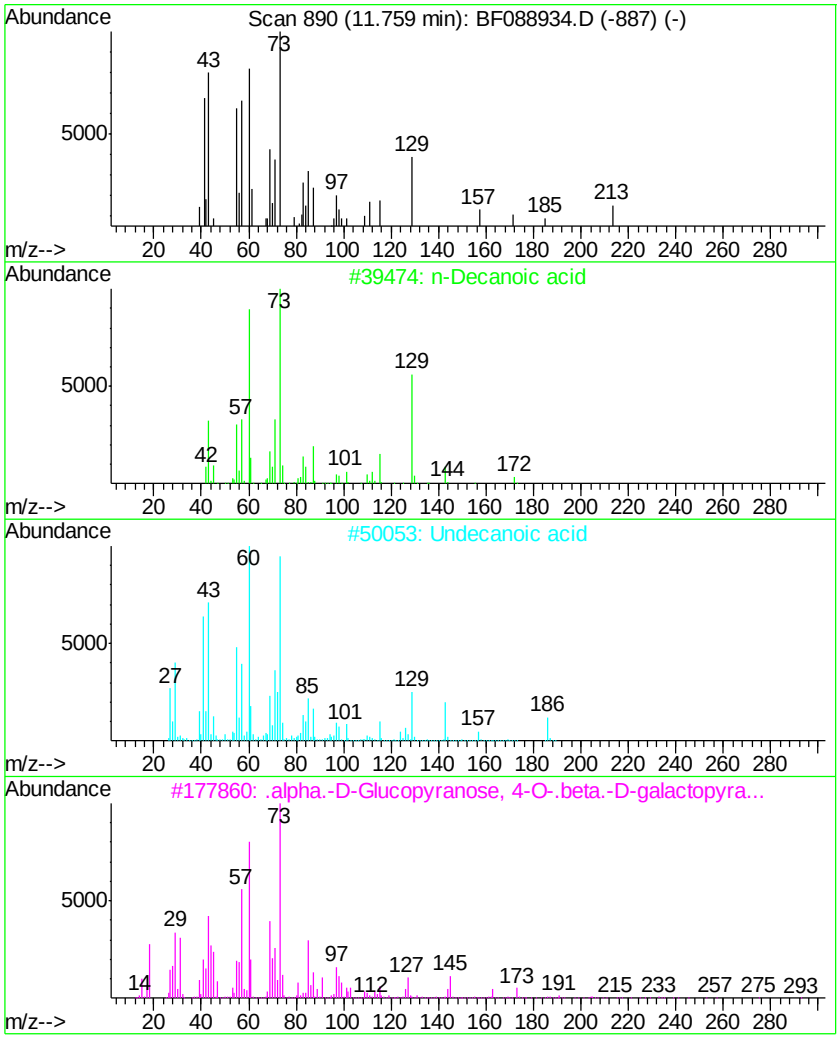
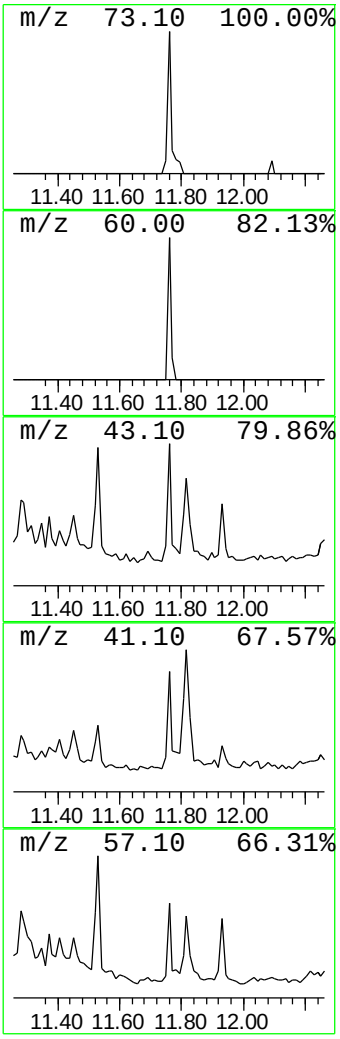
Peak Number 13 n-Decanoic acid Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	2.12 ng	39602	Phenanthrene-d10	11.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	53
2	Undecanoic acid	186	C11H22O2	000112-37-8	50
3	.alpha.-D-Glucopyranose, 4-O-.be...	342	C12H22O11	014641-93-1	50
4	Lactose	342	C12H22O11	000063-42-3	38
5	Octane, 1-(ethenylthio)-	172	C10H20S	042779-08-8	32





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Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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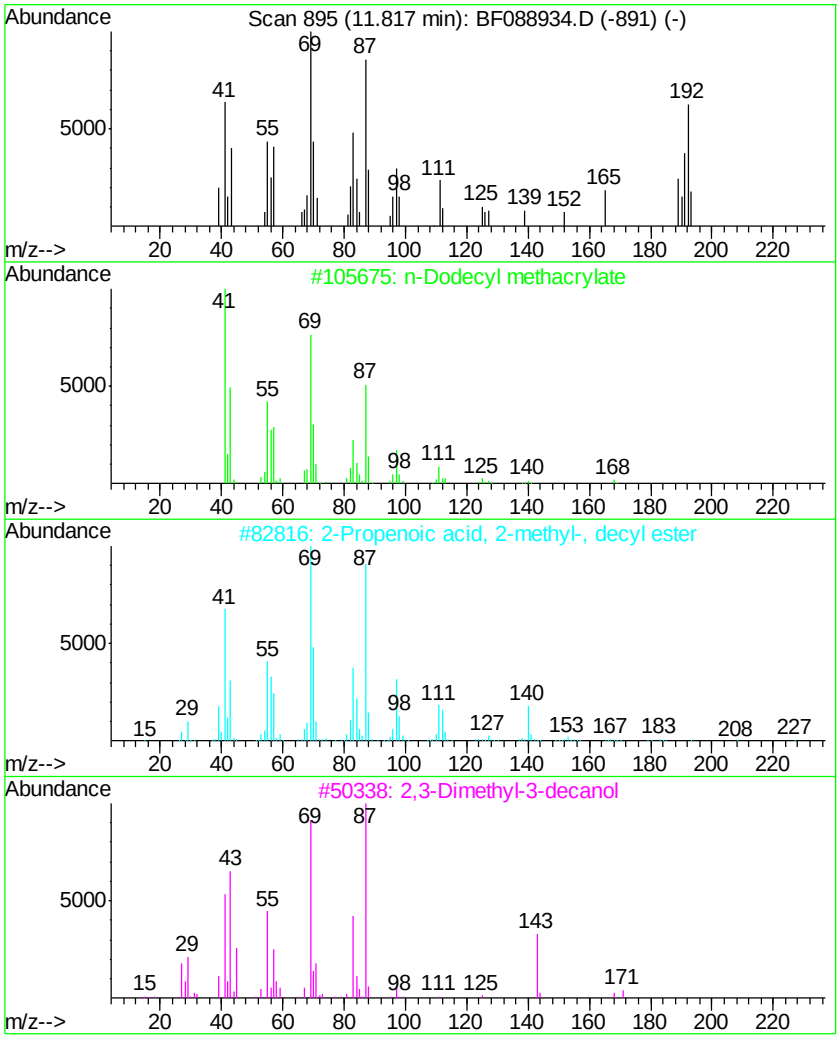
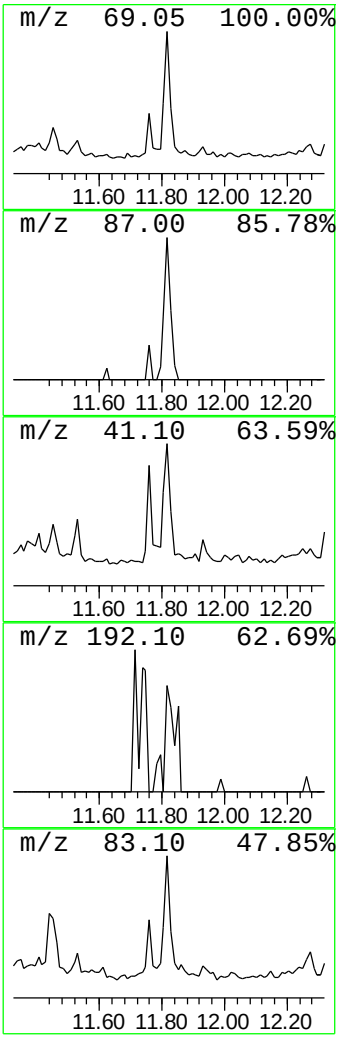
Peak Number 14 n-Dodecyl methacrylate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	4.50 ng	84097	Phenanthrene-d10	11.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	64
2	2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	53
3	2,3-Dimethyl-3-decanol	186	C12H26O	219667-42-2	47
4	Cyclopropanecarboxylic acid,decy...	226	C14H26O2	054460-47-8	43
5	Cyclopropanecarboxylic acid, non...	212	C13H24O2	060128-06-5	43





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Data File : BF088934.D
Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
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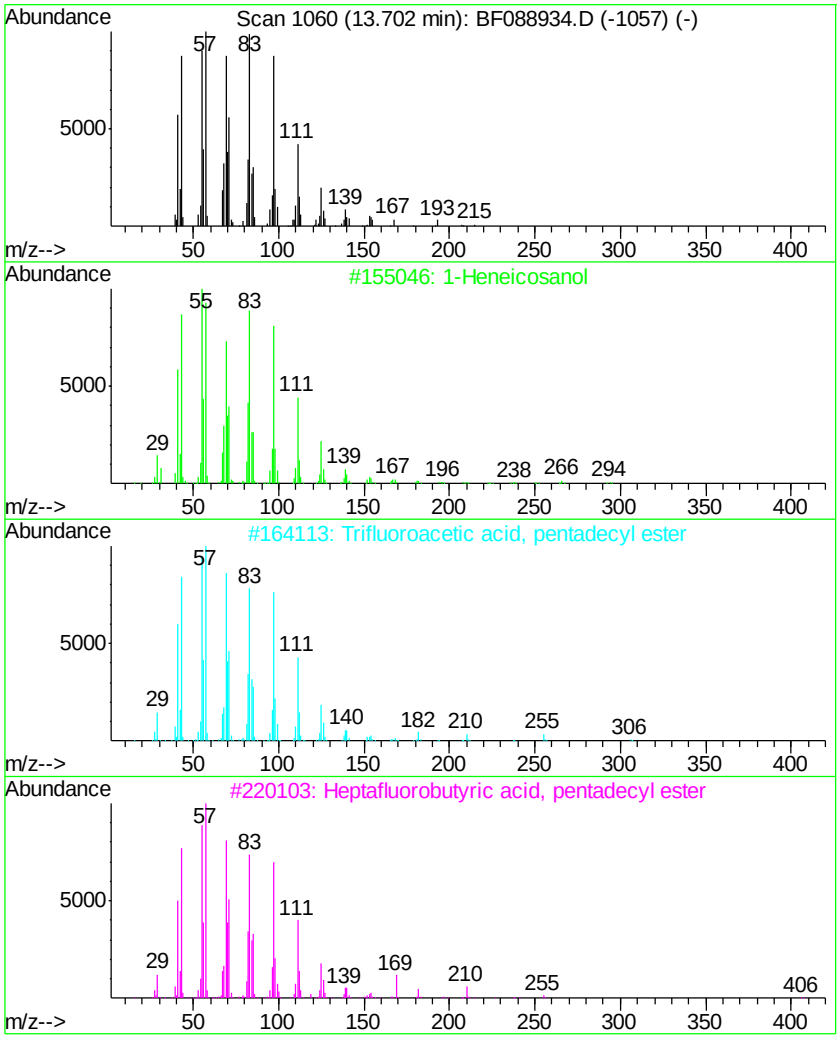
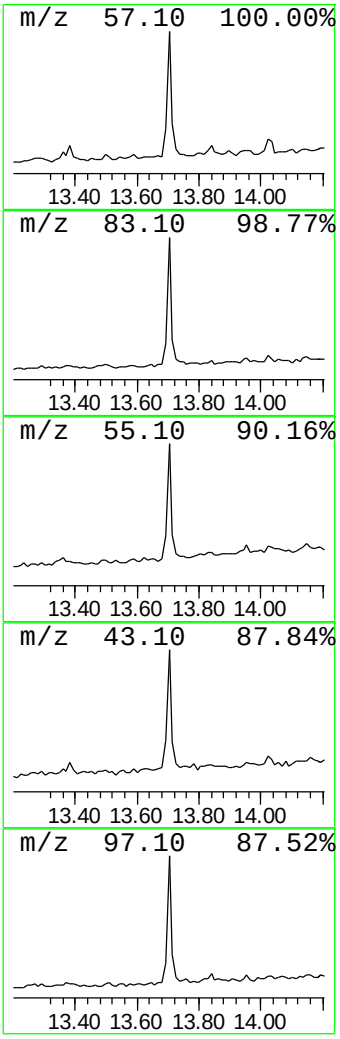
Peak Number 15 1-Heneicosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.70	4.70 ng	123884	Chrysene-d12	13.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	95
3	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	95
4	Behenic alcohol	326	C22H46O	000661-19-8	95
5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	93





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Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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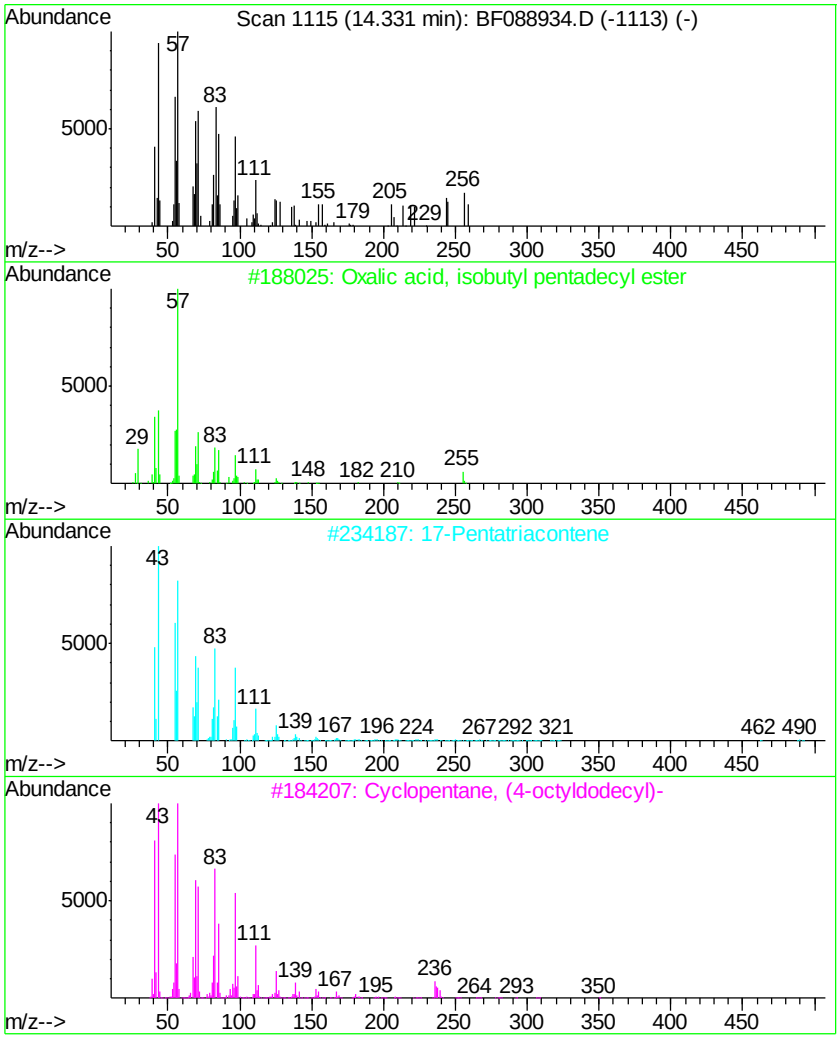
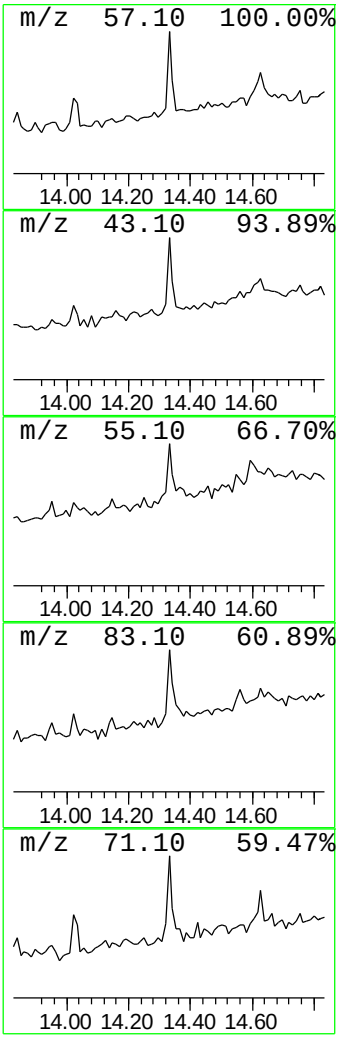
Peak Number 16 Oxalic acid, isobutyl penta... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	2.67 ng	70483	Chrysene-d12	13.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	74
2	17-Pentatriacontene	491	C35H70	006971-40-0	70
3	Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	68
4	Oxalic acid, cyclobutyl octadecy...	396	C24H44O4	1000309-70-8	64
5	Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	58





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Operator : UM/SJ
Sample : H3934-15
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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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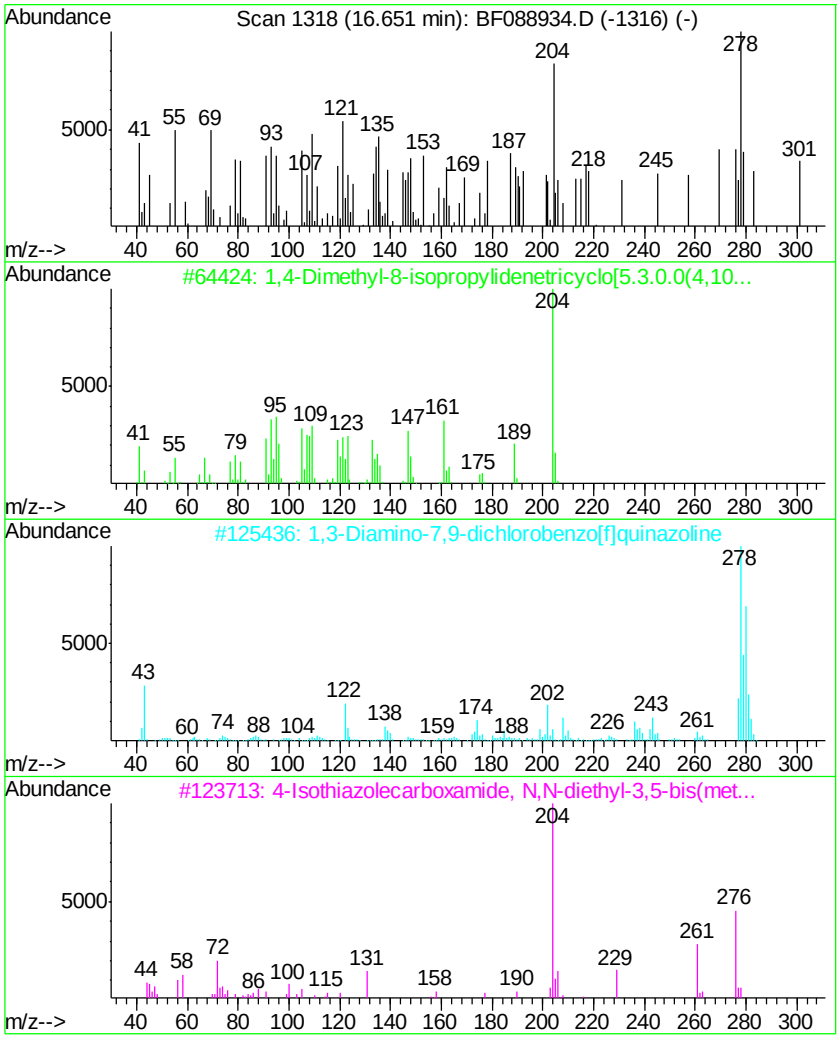
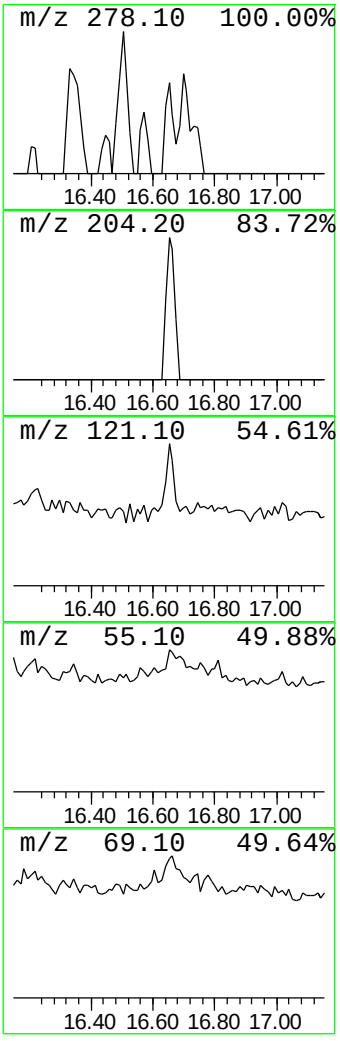
Peak Number 19 unknown16.65 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	2.87 ng	48723	Perylene-d12	15.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	25
2	1,3-Diamino-7,9-dichlorobenzo[f]...	278	C12H8Cl2N4	037521-58-7	25
3	4-Isothiazolecarboxamide, N,N-di...	276	C10H16N2OS3	037572-36-4	22
4	Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	15
5	3,5-Dichloro-2,4-dimethyl-1-meth...	204	C9H10Cl2O	1000250-17-8	15





Tentatively Identified Compound (LSC) summary

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Acq On : 12 Jul 2016 22:27
Operator : UM/SJ
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : C2.3-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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
Propane, 1,1-dime...	1.90	5.4	ng	104968	1	6.71	387943	20.0
unknown2.87	2.87	2.2	ng	42889	1	6.71	387943	20.0
2-Pentanone, 4-hy...	4.86	10.1	ng	195731	1	6.71	387943	20.0
unknown6.48	6.48	91.6	ng	1777460	1	6.71	387943	20.0
Safrole	8.62	5.6	ng	143222	2	7.99	508488	20.0
Juglone	9.86	2.2	ng	50981	3	9.74	457936	20.0
Tridecane	10.65	2.5	ng	46607	4	11.21	374084	20.0
Heptadecane	11.10	2.3	ng	42167	4	11.21	374084	20.0
n-Decanoic acid	11.76	2.1	ng	39602	4	11.21	374084	20.0
n-Dodecyl methacr...	11.82	4.5	ng	84097	4	11.21	374084	20.0
1-Heneicosanol	13.70	4.7	ng	123884	5	13.84	527063	20.0
Oxalic acid, isob...	14.33	2.7	ng	70483	5	13.84	527063	20.0
unknown16.65	16.65	2.9	ng	48723	6	15.21	339182	20.0