LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

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-BF092815.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	2.240	7	9	14	rVB2	11360	27206	1.12%	0.100%
2	5.097	256	259	261	rBV	790409	1134799	46.87%	4.167%
3	5.497	292	294	297	rBV	1623178	1878968	77.61%	6.899%
4	6.526	382	384	387	rBV	1503496	1853130	76.54%	6.804%
5	6.617	390	392	394	rVB	51607	56023	2.31%	0.206%
6	6.675	394	397	399	rBV	1739872	2228382	92.04%	8.182%
7	6.903	415	417	428	rBV	494952	617026	25.48%	2.265%
8	7.063	428	431	433	rBV	1693531	1718151	70.96%	6.308%
9	7.475	464	467	469	rBV	975567	1210250	49.99%	4.444%
10	8.195	527	530	532	rBV	579790	697988	28.83%	2.563%
11	8.240	533	534	536	rVB	44685	57615	2.38%	0.212%
12	8.389	545	547	550	rVB	27697	30301	1.25%	0.111%
13	8.480	550	555	558	rBV4	22364	47856	1.98%	0.176%
14	8.606	564	566	567	rBV	101174	88410	3.65%	0.325%
15	8.720	575	576	580	rBV2	21060	45128	1.86%	0.166%
16	8.835	583	586	588	rVV4	18788	55465	2.29%	0.204%
17	8.869	588	589	591	rVB2	31434	34164	1.41%	0.125%
18	8.972	596	598	601	rBV4	17714	48819	2.02%	0.179%
19	9.098	607	609	611	rVB3	21857	25362	1.05%	0.093%
20	9.212	617	619	621	rVB	114743	148340	6.13%	0.545%
21	9.269	621	624	626	rBV	2494907	2421169	100.00%	8.890%
22	9.349	629	631	633	rVB2	54707	83204	3.44%	0.305%
23	9.623	654	655	656	rBV	19316	25235	1.04%	0.093%
24	9.658	656	658	660	rVV	519812	583411	24.10%	2.142%
25	9.806	669	671	673	rBV3	40653	45002	1.86%	0.165%
26	9.852	673	675	676	rBV2	51098	50785	2.10%	0.186%
27	9.886	676	678	679	rVB	20483	27661	1.14%	0.102%
28	9.943	681	683	685	rVB	829328	819777	33.86%	3.010%
29	10.046	691	692	694	rBV2	25927	28343	1.17%	0.104%
30	10.138	698	700	703	rVB3	43991	90526	3.74%	0.332%
31	10.195	703	705	709	rVB3	69328	99585	4.11%	0.366%
32	10.263	709	711	713	rBV2	44004	63259	2.61%	0.232%
33	10.309	713	715	717	rVB2	37169	53700	2.22%	0.197%
34	10.366	717	720	722	rBV2	44342	116712	4.82%	0.429%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

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-BF092815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.435	724	726	728	rVB3	29503	41383	1.71%	0.152%
36	10.492	728	731	732	rBV3	29009	60779	2.51%	0.223%
37	10.595	737	740	743	rVB	211071	272567	11.26%	1.001%
38	10.641	743	744	747	rVB3	24735	28543	1.18%	0.105%
39	10.732	750	752	755	rBV	1280519	1401011	57.87%	5.144%
40	10.858	760	763	765	rVV	495109	526669	21.75%	1.934%
41	10.926	767	769	773	rBV5	32266	77738	3.21%	0.285%
42	11.304	799	802	805	rBV2	1453117	2199109	90.83%	8.074%
43	11.361	805	807	809	rVB	690541	738224	30.49%	2.710%
44	11.406	809	811	812	rBV	126075	109223	4.51%	0.401%
45	11.429	812	813	821	rVB	716367	904514	37.36%	3.321%
46	11.669	832	834	836	rBV	73127	86005	3.55%	0.316%
47	11.715	836	838	840	rVV2	29220	38277	1.58%	0.141%
48	11.761	840	842	846	rVB4	26864	56824	2.35%	0.209%
49	12.412	897	899	903	rVB3	18827	28442	1.17%	0.104%
50	12.481	903	905	907	rBV2	21116	27719	1.14%	0.102%
51	12.641	918	919	926	rVB	31185	56914	2.35%	0.209%
52	12.869	937	939	943	rBV	27907	50339	2.08%	0.185%
53	13.018	950	952	964	rVB	2230975	2351374	97.12%	8.633%
54	13.910	1028	1030	1040	rBV	86479	226265	9.35%	0.831%
55	14.070	1042	1044	1057	rVB	592492	842237	34.79%	3.092%
56	15.110	1132	1135	1139	rBV2	9099	25676	1.06%	0.094%
57	15.521	1169	1171	1189	rBV	302280	574573	23.73%	2.110%

Sum of corrected areas: 27236157

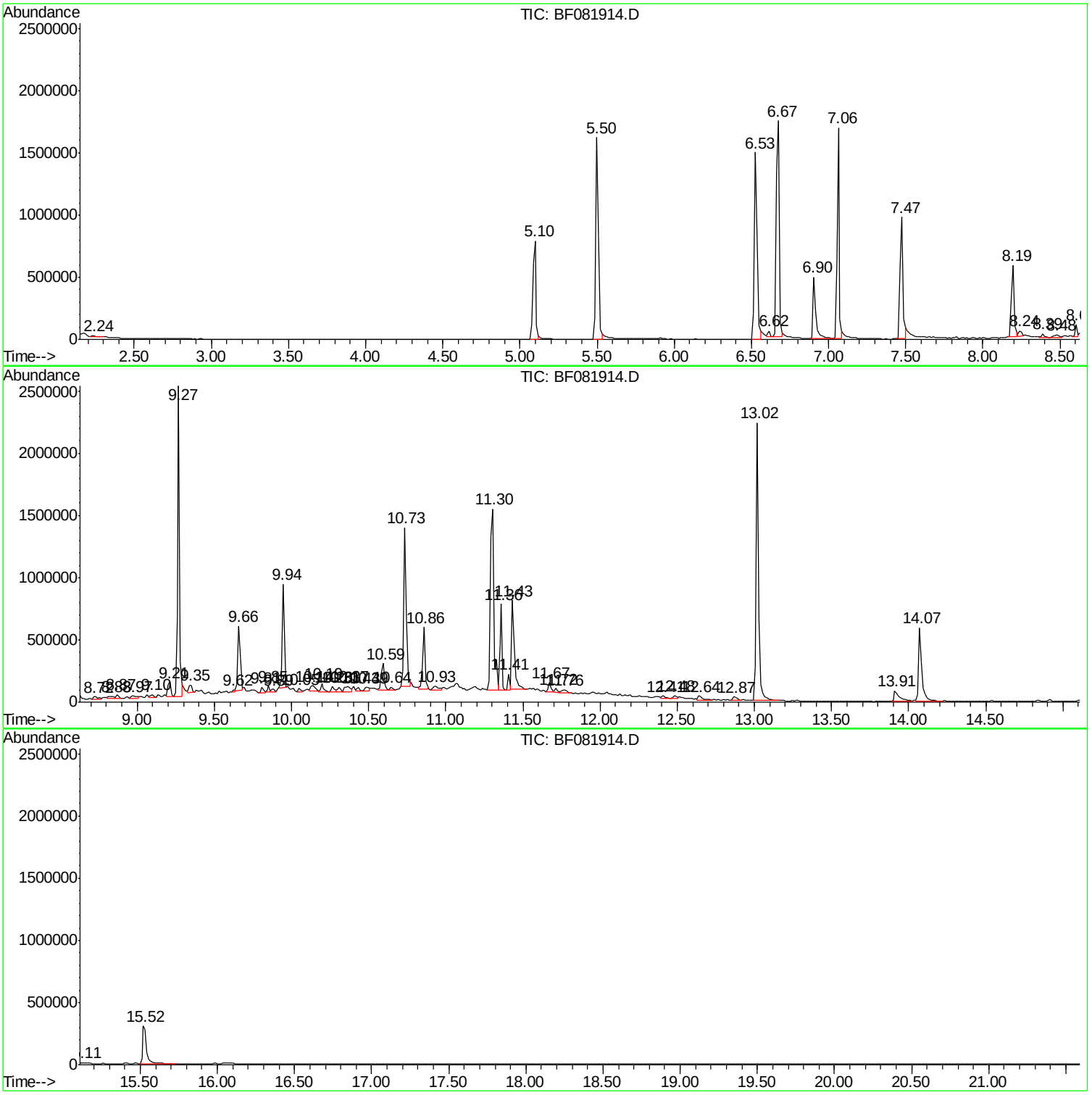
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

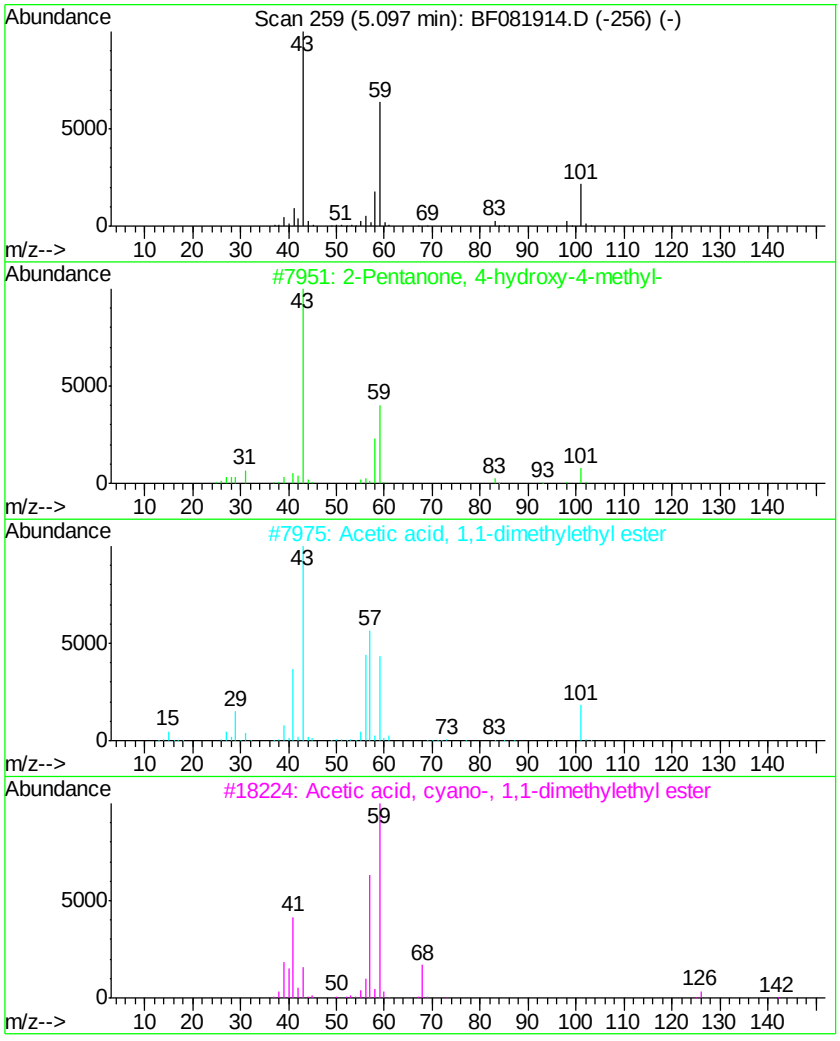
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

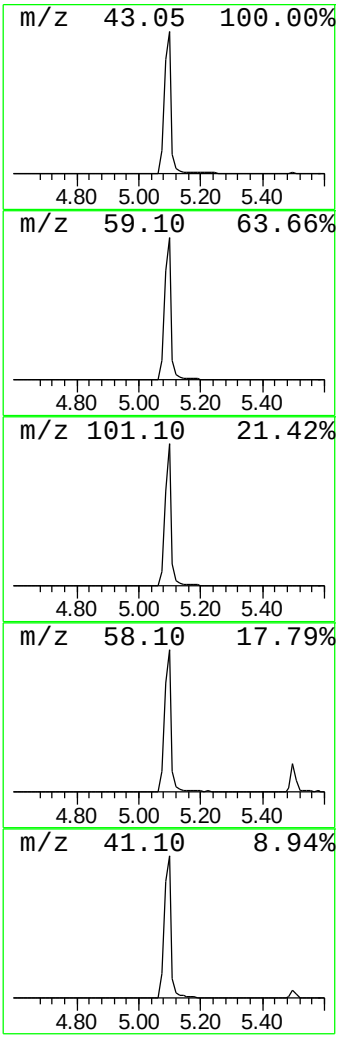
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	36.78 ng	1134800	1,4-Dichlorobenzene-d4	6.90

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, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	Acetone	58	C3H6O	000067-64-1	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

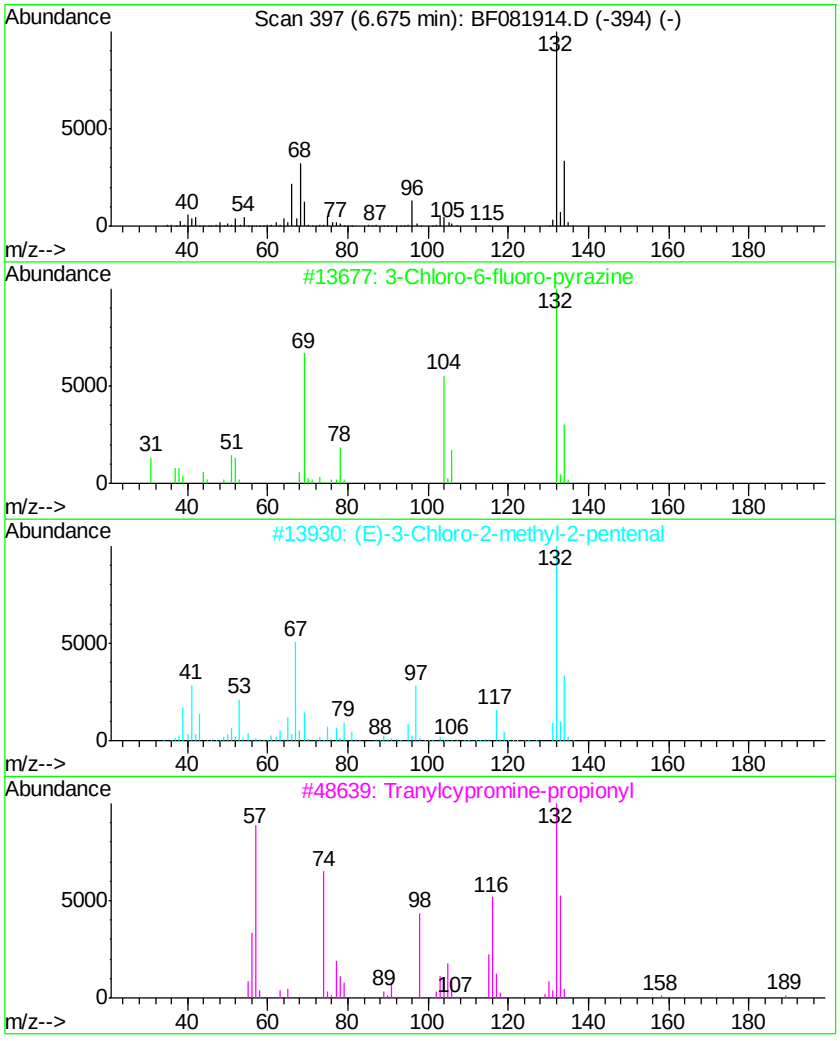
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

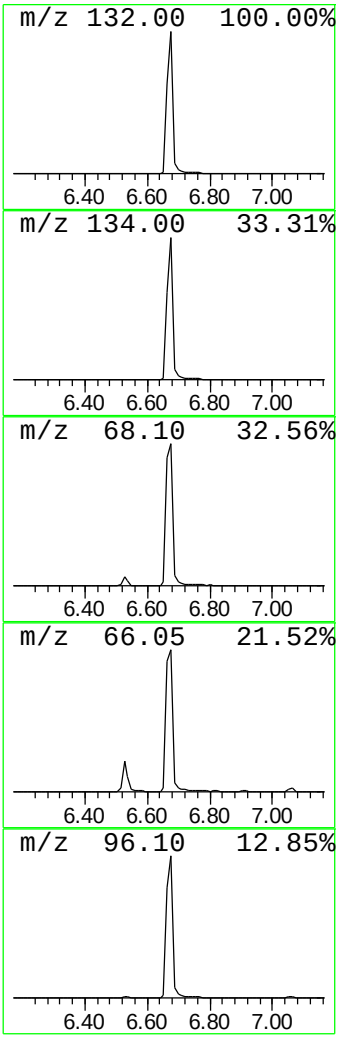
****

Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	72.23 ng	2228380	1,4-Dichlorobenzene-d4	6.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	Amphetaminil	250	C17H18N2	017590-01-1	12
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

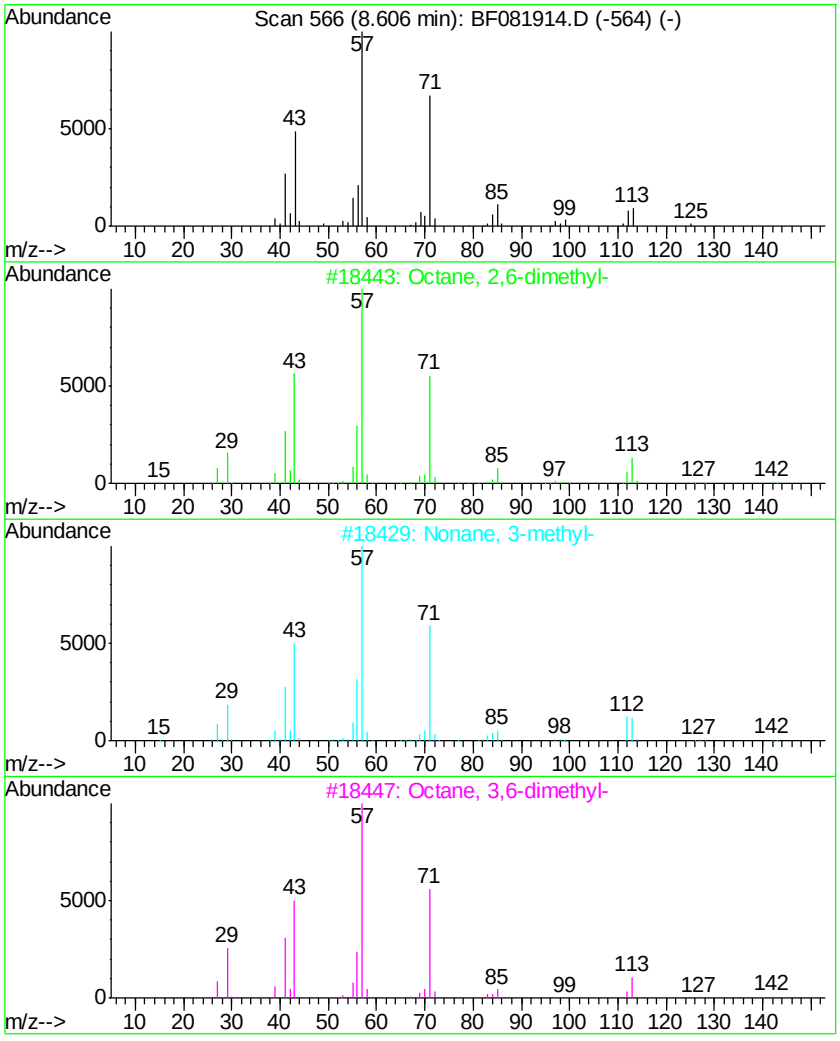
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

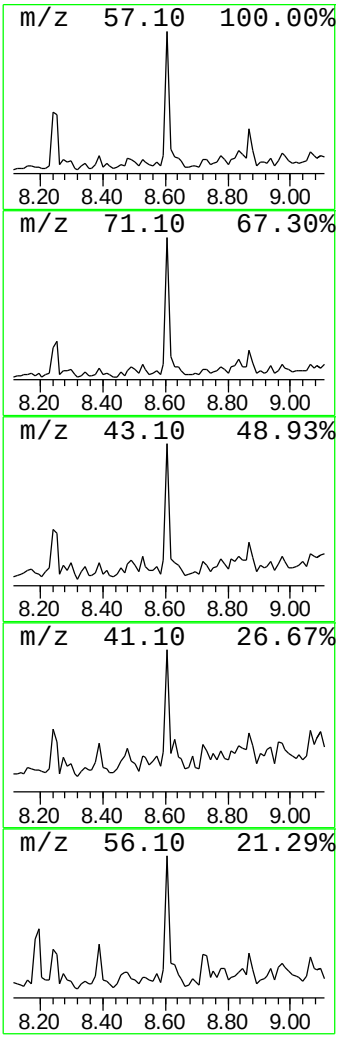
****

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	2.53 ng	88410	Naphthalene-d8	8.19

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	78
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	78
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
4	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	72
5	Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	72





Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

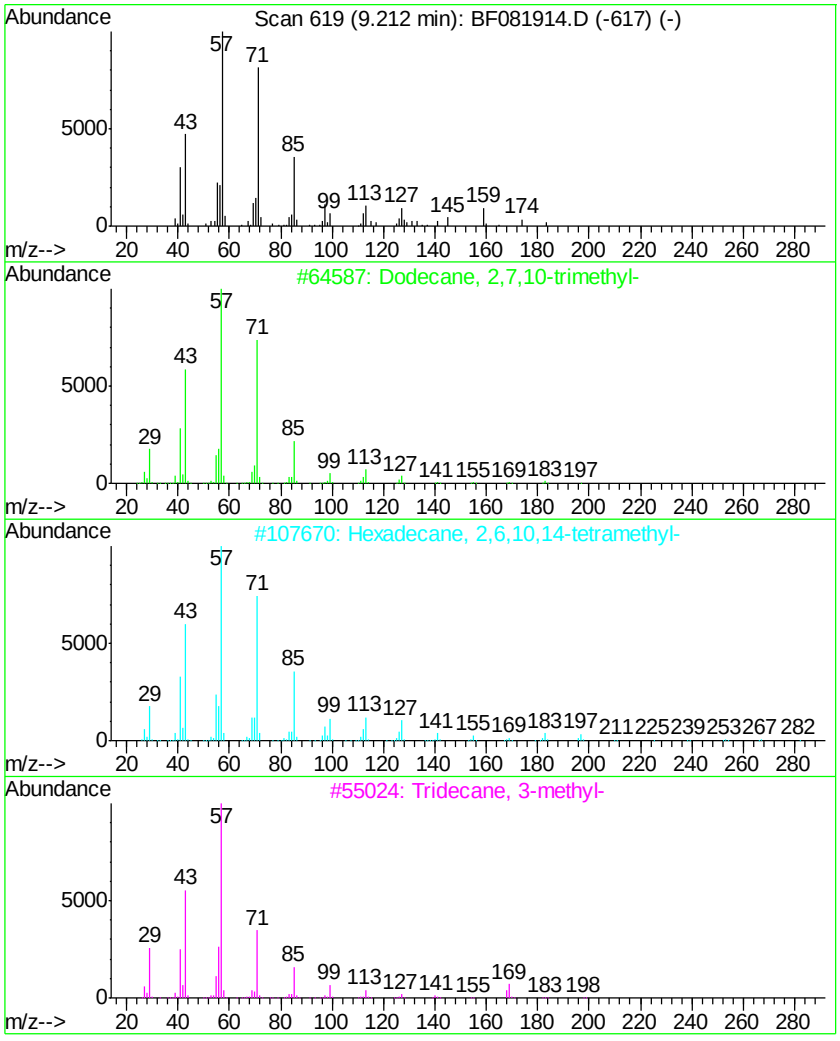
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

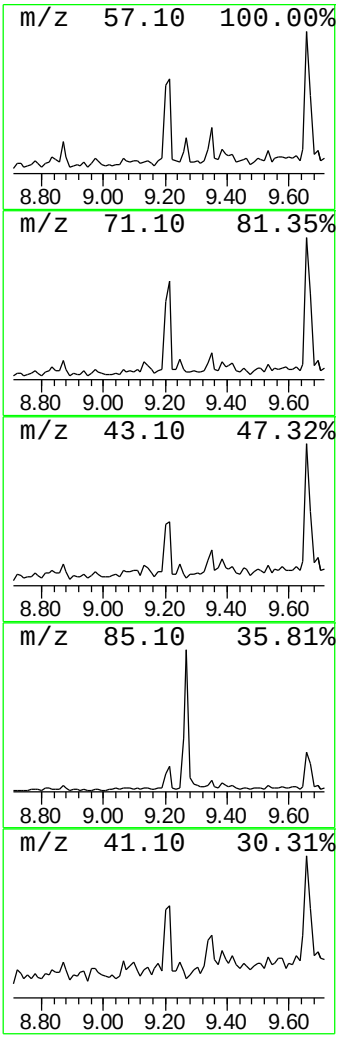
****

Peak Number 5 Dodecane, 2,7,10-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	3.62 ng	148340	Acenaphthene-d10	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78
3	Tridecane, 3-methyl-	198	C14H30	006418-41-3	72
4	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
5	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	64





Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

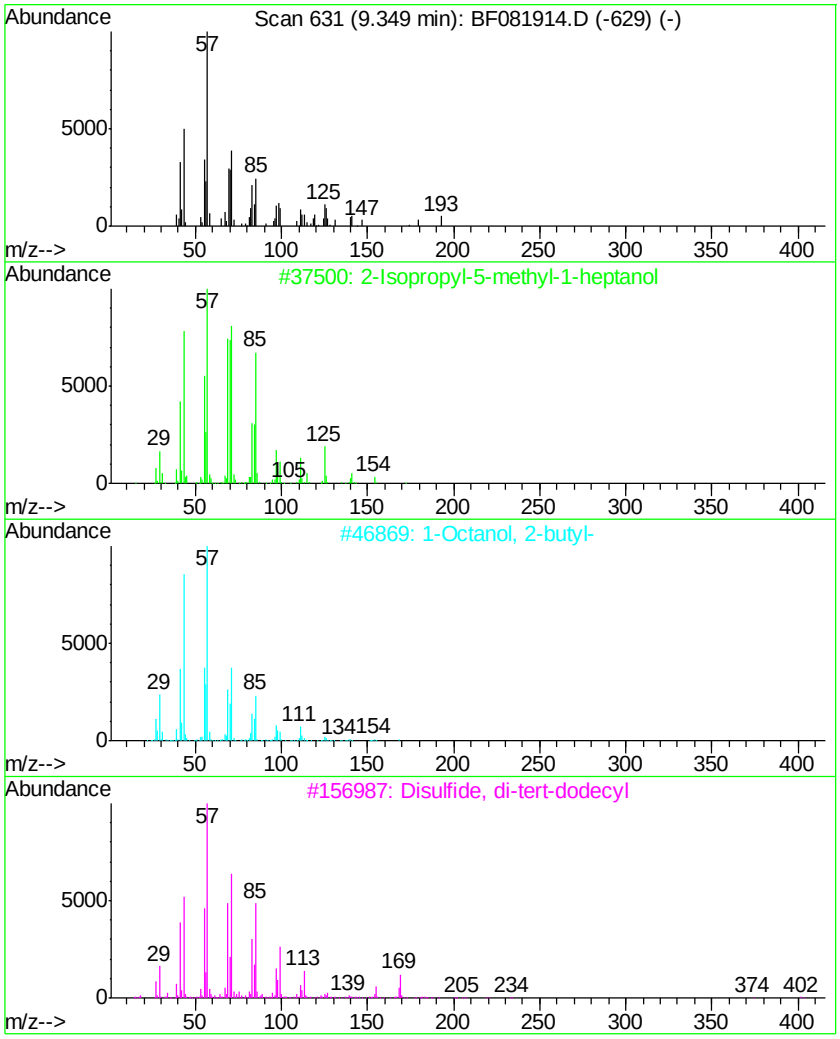
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

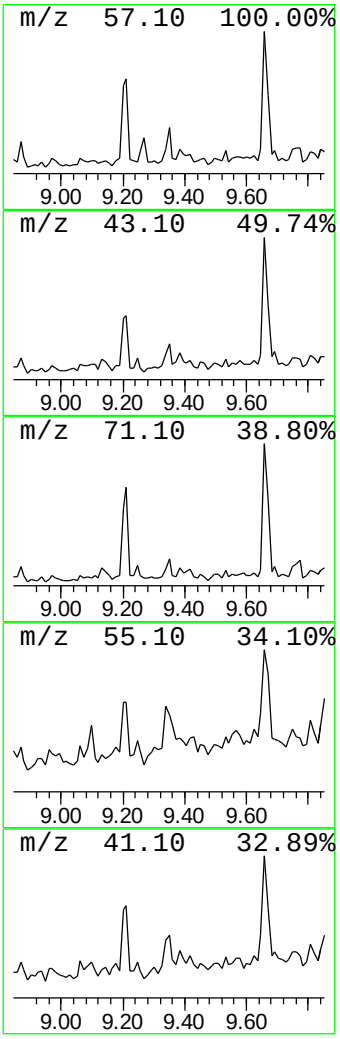
****

Peak Number 6 2-Isopropyl-5-methyl-1-hept... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.35	2.03 ng	83204	Acenaphthene-d10	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Isopropyl-5-methyl-1-heptanol	172	C11H24O	091337-07-4	59
2	1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	59
3	Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	53
4	Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	53
5	Tridecane, 6-methyl-	198	C14H30	013287-21-3	50





Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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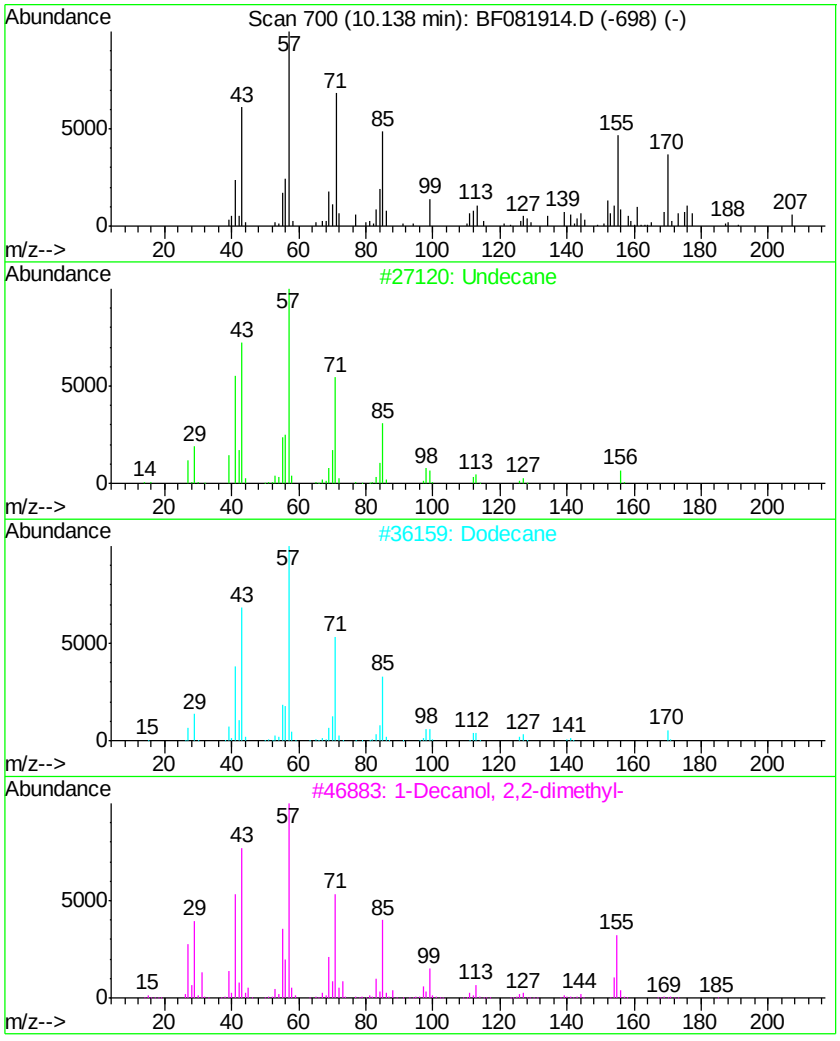
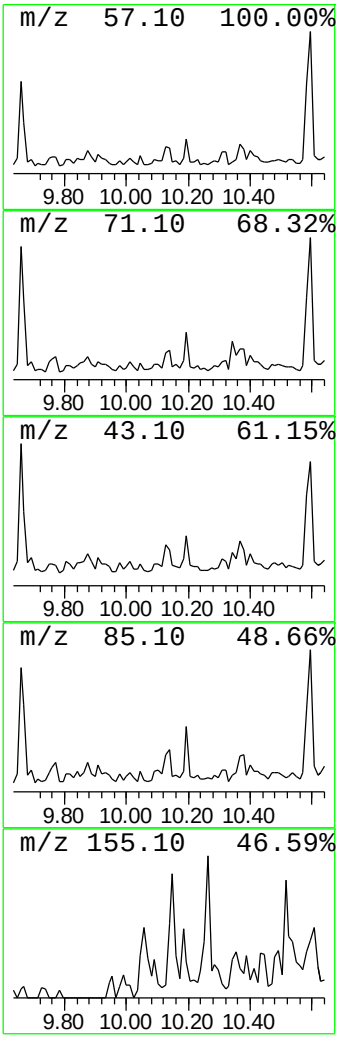
Peak Number 7 Undecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.21 ng	90526	Acenaphthene-d10	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	50
2	Dodecane	170	C12H26	000112-40-3	43
3	1-Decanol, 2,2-dimethyl-	186	C12H26O	002370-15-2	43
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	43
5	Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

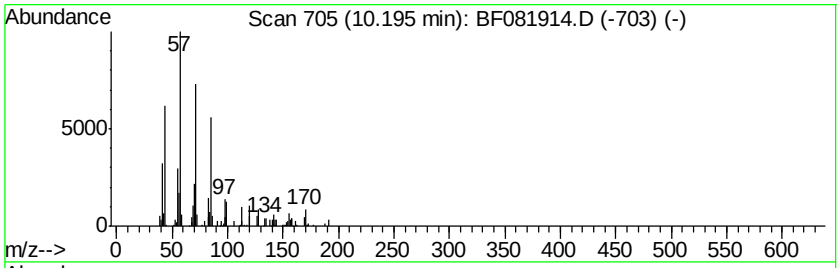
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

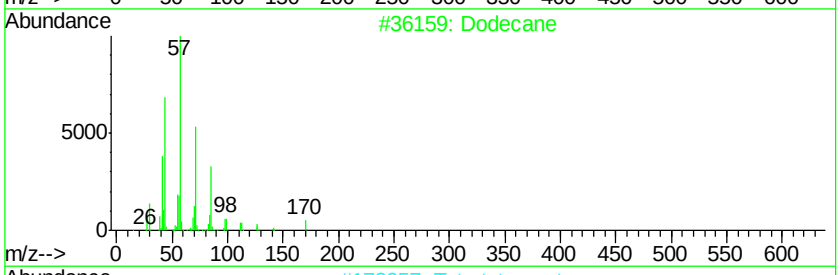
****

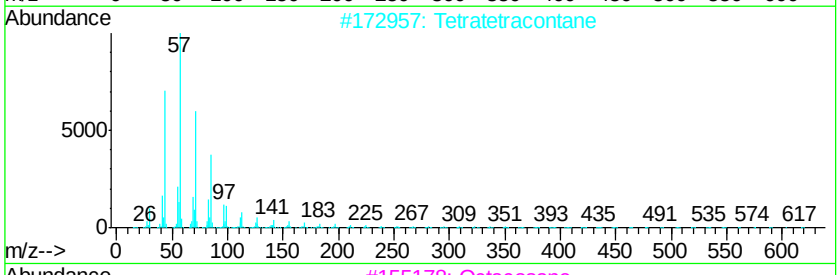
Peak Number 8 Dodecane Concentration Rank 12

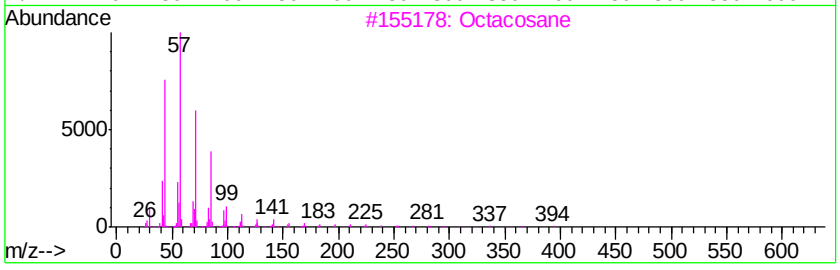
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.19	2.43 ng	99585	Acenaphthene-d10	9.94

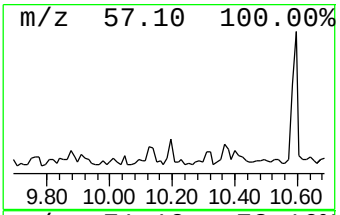
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	87
2	Tetratetracontane	619	C44H90	007098-22-8	83
3	Octacosane	394	C28H58	000630-02-4	80
4	Nonacosane	408	C29H60	000630-03-5	72
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72

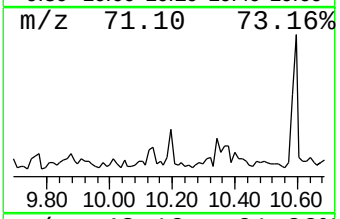


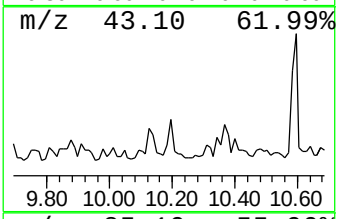


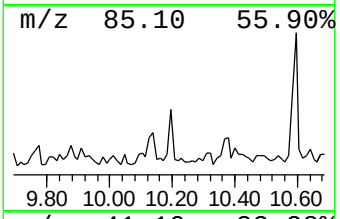


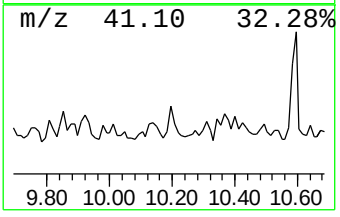












Data Path : Z:\HPCHEM1\BNA F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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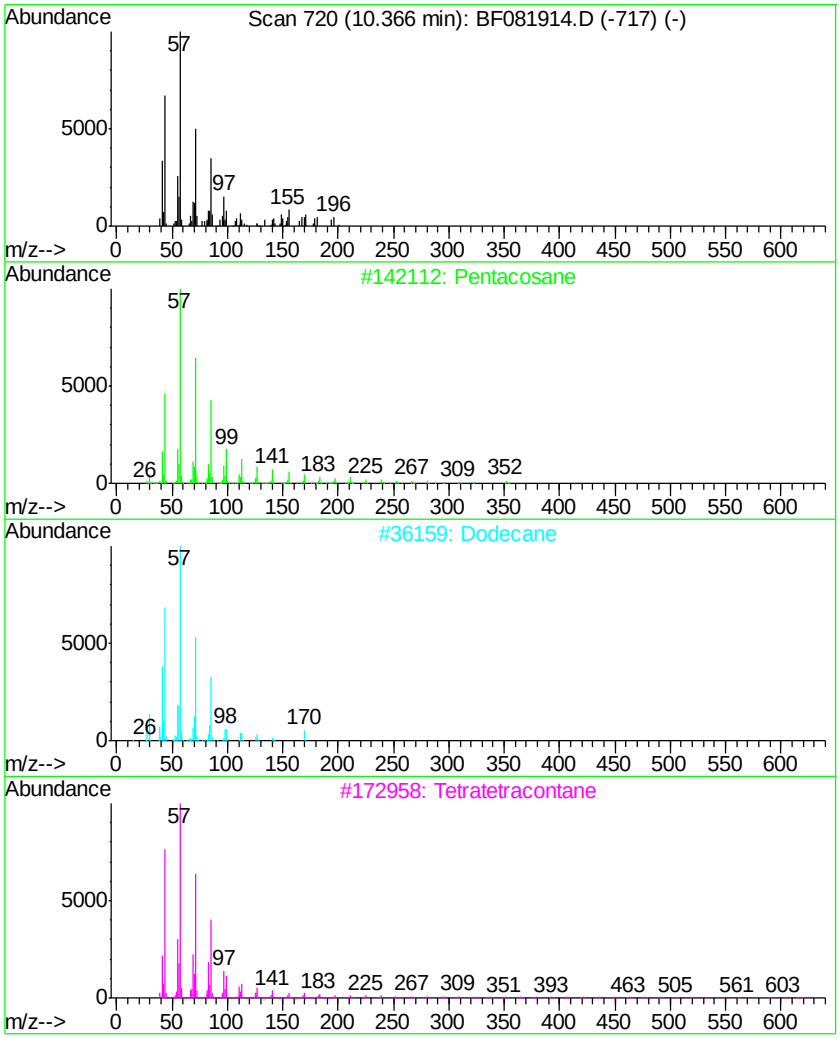
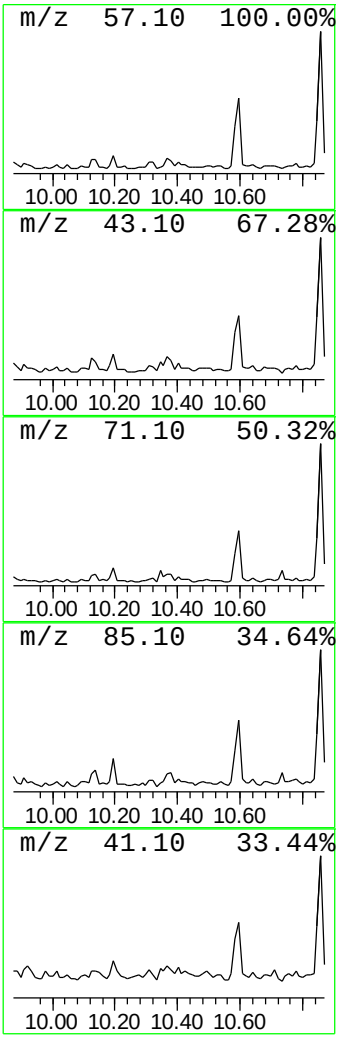
Peak Number 9 Pentacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	2.85 ng	116712	Acenaphthene-d10	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	64
2	Dodecane	170	C12H26	000112-40-3	62
3	Tetratetracontane	619	C44H90	007098-22-8	59
4	Hexatriacontane	507	C36H74	000630-06-8	59
5	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	53





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Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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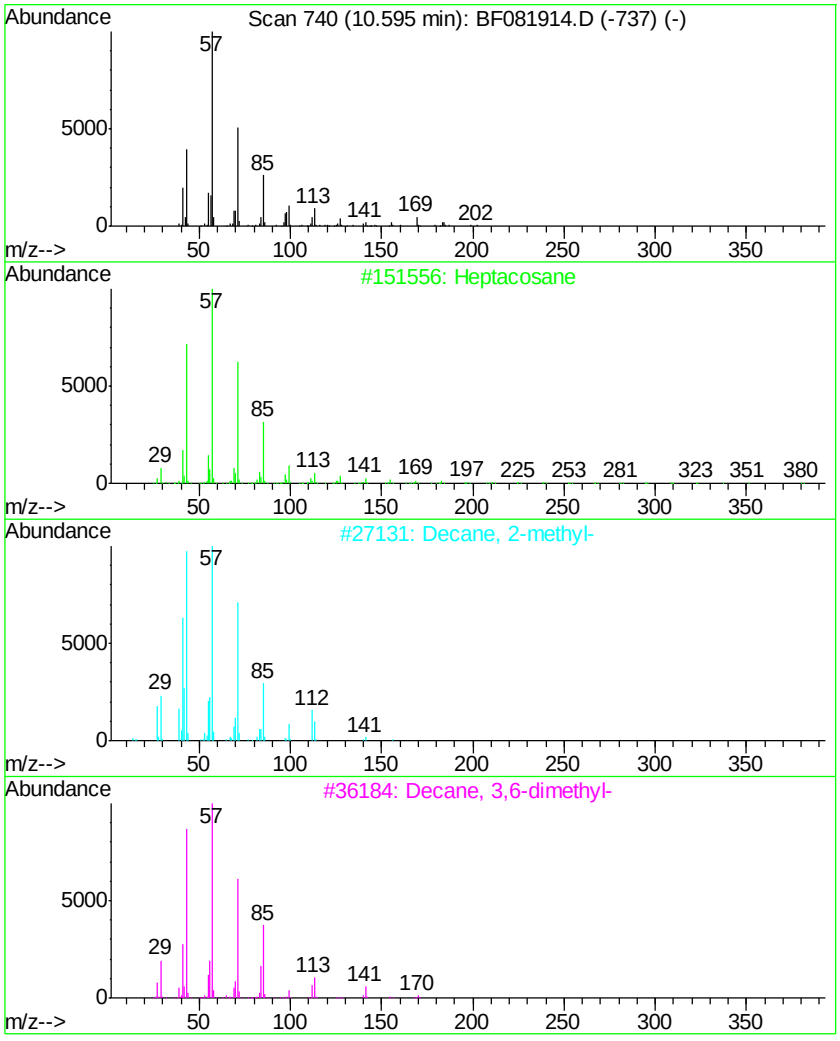
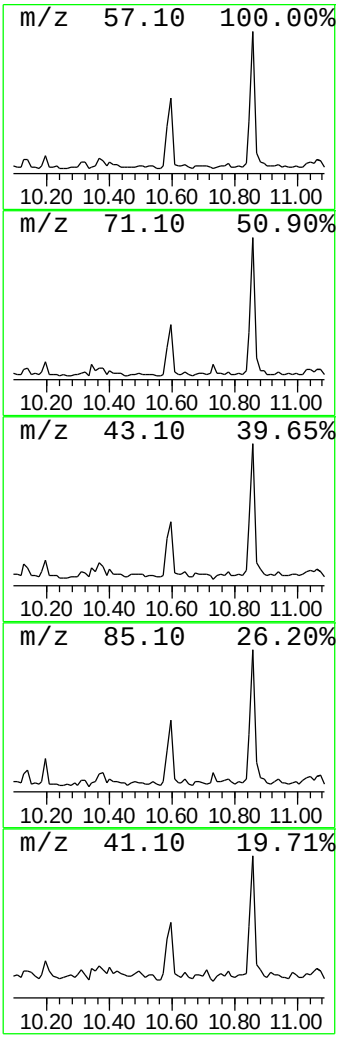
Peak Number 10 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.59	6.65 ng	272567	Acenaphthene-d10	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	86
2	Decane, 2-methyl-	156	C11H24	006975-98-0	81
3	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	81
4	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	76
5	Tridecane	184	C13H28	000629-50-5	72





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Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
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TIC Integration Parameters: LSCINT.P

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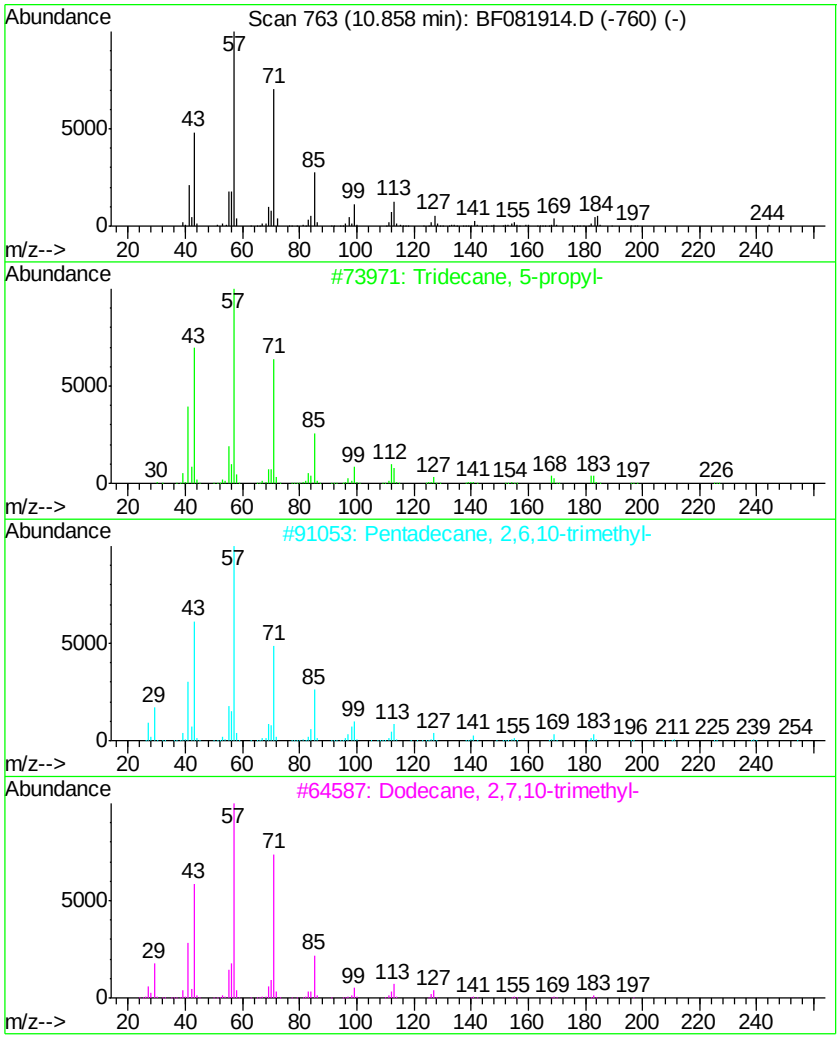
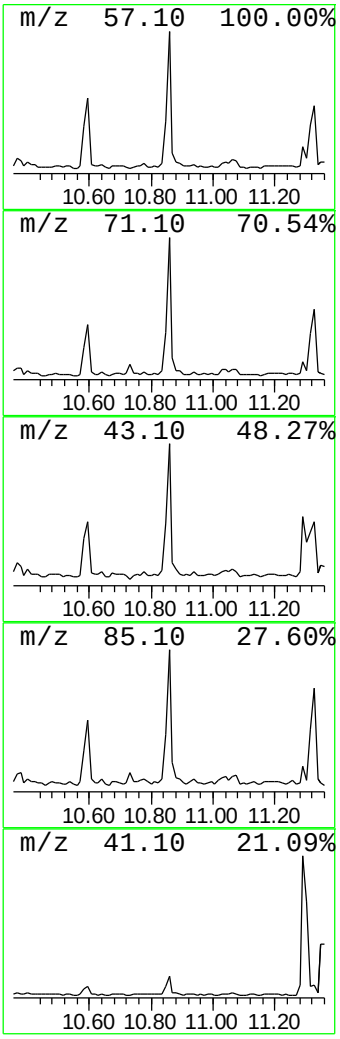
Peak Number 11 Tridecane, 5-propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.86	11.65 ng	526669	Phenanthrene-d10	11.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 5-propyl-	226	C16H34	055045-11-9	86
2	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
3	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
4	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64
5	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	64





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Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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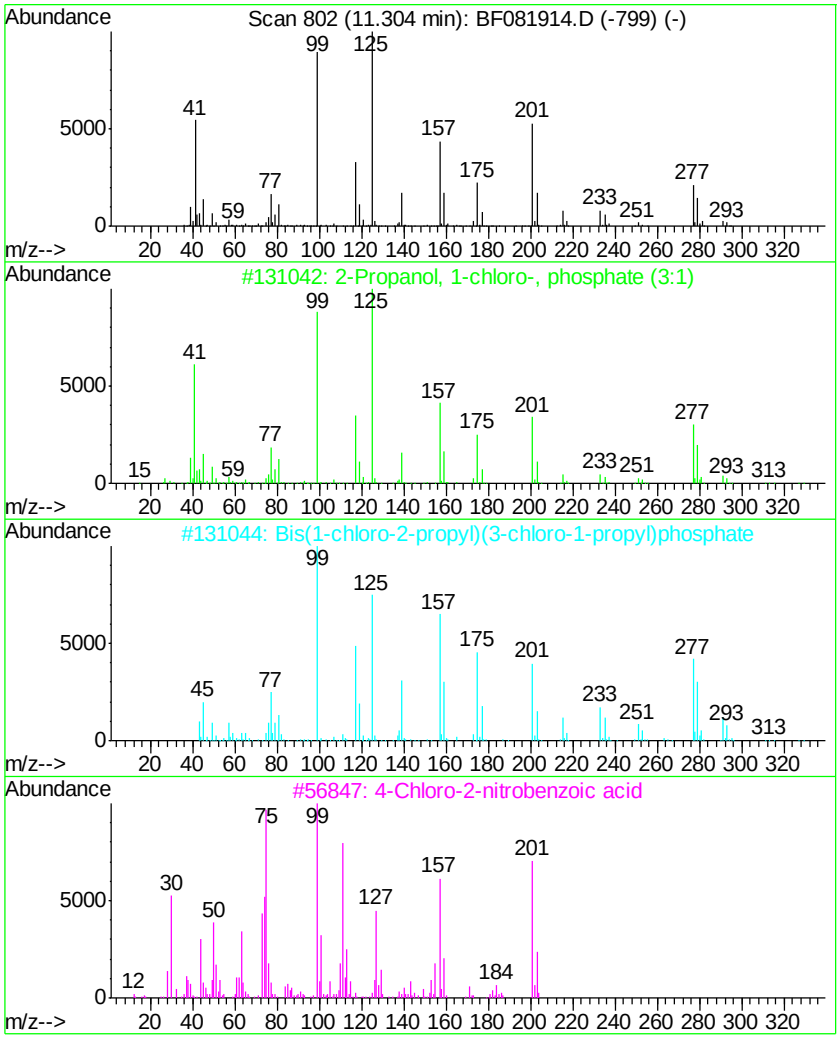
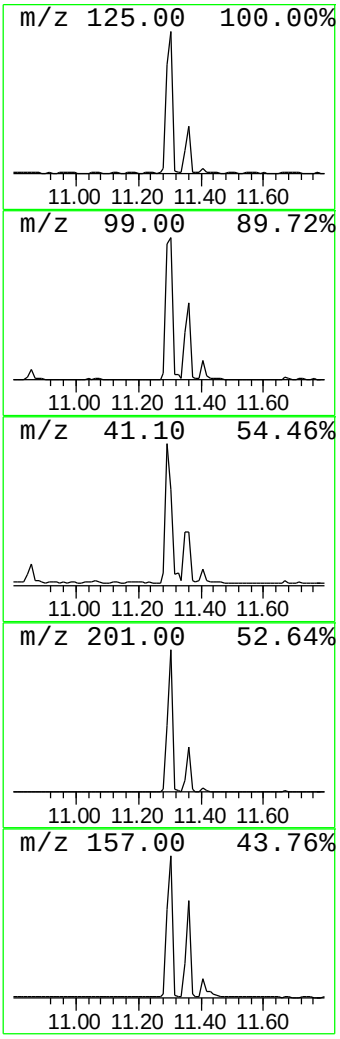
Peak Number 12 2-Propanol, 1-chloro-, phos... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	48.63 ng	2199110	Phenanthrene-d10	11.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	90
2	Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	49
3	4-Chloro-2-nitrobenzoic acid	201	C7H4ClNO4	006280-88-2	22
4	n-pentyl(1-propenyl)dimethylsilane	170	C10H22Si	136964-13-1	10
5	1,2,5-Oxadiazol-3-amine, 4-[5-(4...	277	C11H8ClN5O2	1000276-59-0	9





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Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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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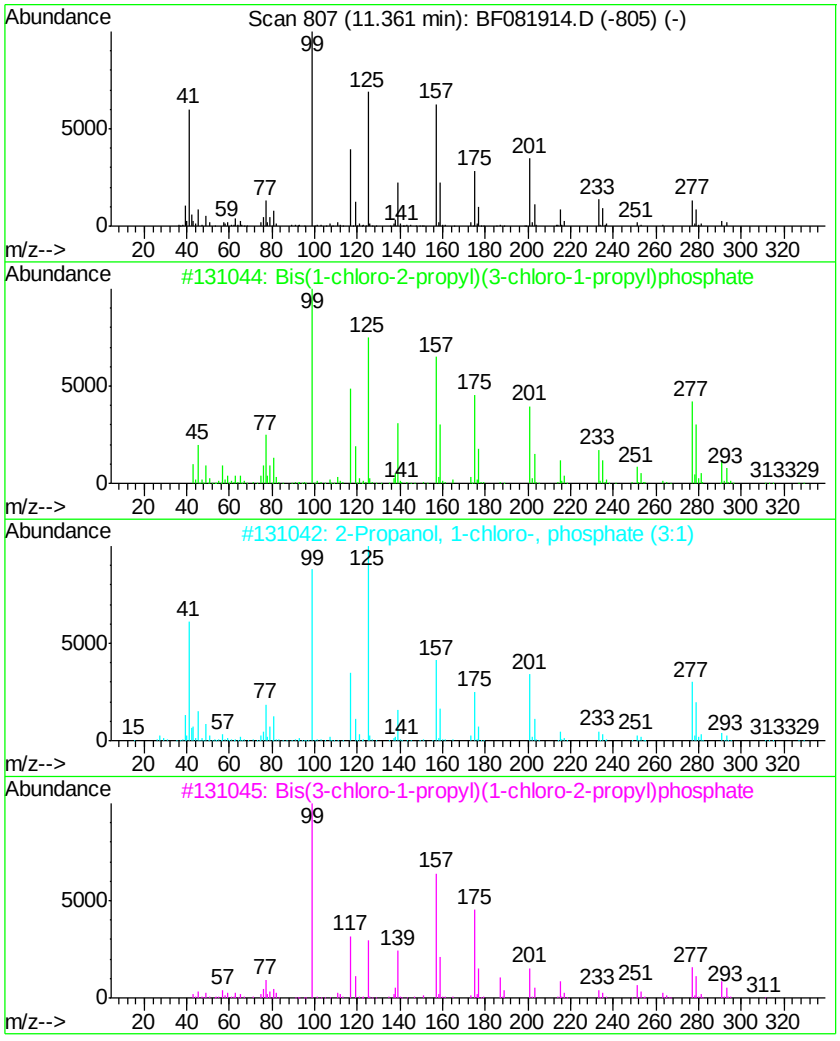
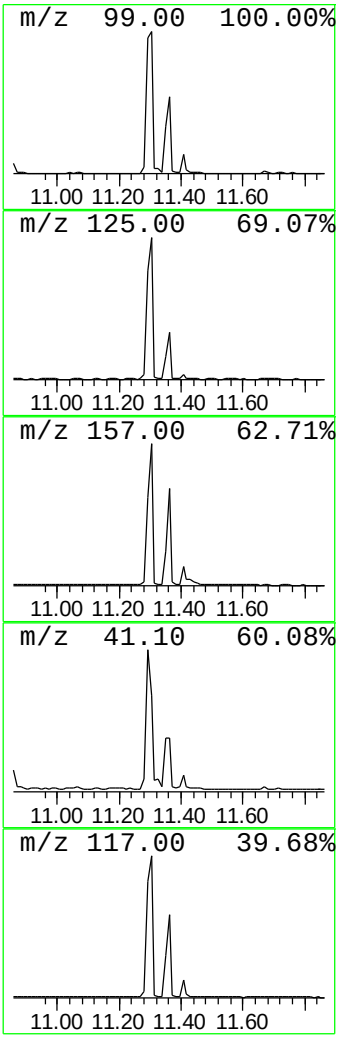
Peak Number 13 Bis(1-chloro-2-propyl)(3-ch... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	16.32 ng	738224	Phenanthrene-d10	11.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	87
2	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	47
3	Bis(3-chloro-1-propyl)(1-chloro-...	326	C9H18Cl3O4P	137888-35-8	42
4	Tris(3-chloropropyl) phosphate	326	C9H18Cl3O4P	001067-98-7	36
5	3-Pyridinecarboxylic acid, 2-chl...	157	C6H4ClNO2	002942-59-8	9





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Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
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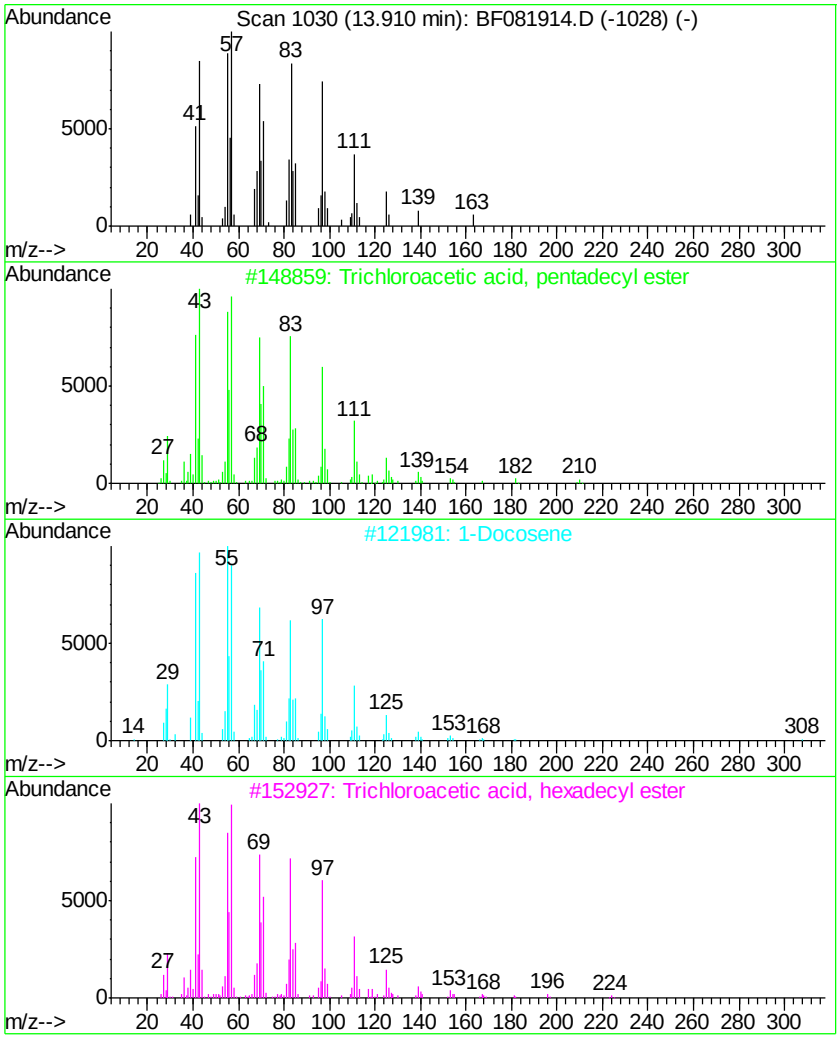
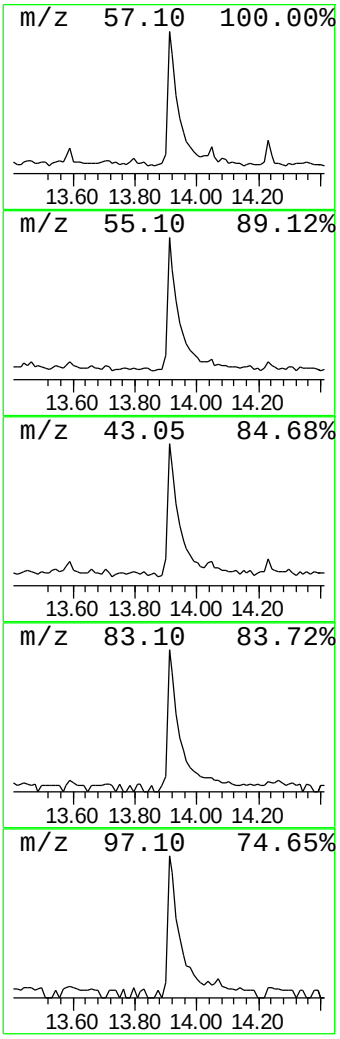
Peak Number 14 Trichloroacetic acid, penta... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	5.37 ng	226265	Chrysene-d12	14.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
2	1-Docosene	308	C22H44	001599-67-3	91
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4	5-Eicosene, (E)-	280	C20H40	074685-30-6	91
5	Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093015\
Data File : BF081914.D
Acq On : 30 Sep 2015 21:19
Operator : UM/IZ
Sample : G3868-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-02(EAST)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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...	5.10	36.8	ng	1134800	1	6.90	617026	20.0
unknown6.67	6.67	72.2	ng	2228380	1	6.90	617026	20.0
Octane, 2,6-dimet...	8.61	2.5	ng	88410	2	8.19	697988	20.0
Dodecane, 2,7,10-...	9.21	3.6	ng	148340	3	9.94	819777	20.0
2-Isopropyl-5-met...	9.35	2.0	ng	83204	3	9.94	819777	20.0
Undecane	10.14	2.2	ng	90526	3	9.94	819777	20.0
Dodecane	10.19	2.4	ng	99585	3	9.94	819777	20.0
Pentacosane	10.37	2.9	ng	116712	3	9.94	819777	20.0
Heptacosane	10.59	6.7	ng	272567	3	9.94	819777	20.0
Tridecane, 5-propyl-	10.86	11.7	ng	526669	4	11.43	904514	20.0
2-Propanol, 1-chl...	11.30	48.6	ng	2199110	4	11.43	904514	20.0
Bis(1-chloro-2-pr...	11.36	16.3	ng	738224	4	11.43	904514	20.0
Trichloroacetic a...	13.91	5.4	ng	226265	5	14.07	842237	20.0