LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

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_G\METHODS\8270-BG022316.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	3.055	32	37	53	rBB	172881	238641	98.10%	11.246%
2	3.196	55	61	68	rBB2	4876	8408	3.46%	0.396%
3	4.659	307	310	315	rBB	2619	3564	1.47%	0.168%
4	5.281	410	416	426	rBB	33663	51127	21.02%	2.409%
5	5.781	495	501	509	rBB	94802	147227	60.52%	6.938%
6	7.408	771	778	787	rBB	110731	182037	74.83%	8.578%
7	7.754	830	837	844	rBB	107951	181456	74.59%	8.551%
8	8.201	908	913	919	rBB	17861	28512	11.72%	1.344%
9	8.536	963	970	976	rBB	77365	130547	53.67%	6.152%
10	9.387	1108	1115	1123	rBB	69755	121369	49.89%	5.719%
11	11.026	1387	1394	1401	rBB	28120	50930	20.94%	2.400%
12	13.447	1800	1806	1813	rBB	166400	243262	100.00%	11.464%
13	14.269	1939	1946	1952	rBB	23806	32850	13.50%	1.548%
14	14.663	2010	2013	2018	rBB	6311	7508	3.09%	0.354%
15	14.821	2034	2040	2046	rBB2	41019	60745	24.97%	2.863%
16	15.608	2170	2174	2179	rVB	7676	9880	4.06%	0.466%
17	16.014	2237	2243	2248	rBB4	2110	4527	1.86%	0.213%
18	16.308	2288	2293	2300	rVB	115638	168783	69.38%	7.954%
19	16.466	2314	2320	2329	rBV	7503	13356	5.49%	0.629%
20	16.537	2329	2332	2338	rVB2	2577	3352	1.38%	0.158%
21	16.631	2341	2348	2353	rBV	6819	9771	4.02%	0.460%
22	16.689	2353	2358	2364	rVB3	5932	10311	4.24%	0.486%
23	16.748	2364	2368	2372	rBV2	6402	8234	3.38%	0.388%
24	16.795	2372	2376	2381	rVV3	3153	4769	1.96%	0.225%
25	16.954	2396	2403	2409	rVB6	4000	9178	3.77%	0.433%
26	17.012	2409	2413	2417	rBV	6635	9430	3.88%	0.444%
27	17.083	2422	2425	2429	rVB3	2821	4009	1.65%	0.189%
28	17.259	2451	2455	2460	rBV	3427	4779	1.96%	0.225%
29	17.559	2501	2506	2511	rBV	38420	53809	22.12%	2.536%
30	17.600	2511	2513	2519	rVB	3820	4300	1.77%	0.203%
31	18.411	2648	2651	2658	rBV2	5693	8376	3.44%	0.395%
32	19.591	2849	2852	2857	rBV	4715	5969	2.45%	0.281%
33	19.967	2910	2916	2919	rVB3	7656	11459	4.71%	0.540%
34	20.138	2940	2945	2950	rBV	142939	180055	74.02%	8.485%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

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_G\METHODS\8270-BG022316.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	20.813	3057	3060	3066	rVB2	7470	9077	3.73%	0.428%
36	21.412	3157	3162	3168	rBV2	10756	16492	6.78%	0.777%
37	21.829	3227	3233	3239	rBV	27210	42150	17.33%	1.986%
38	24.091	3613	3618	3621	rBV6	2692	6085	2.50%	0.287%
39	25.154	3790	3799	3810	rBV4	10280	32612	13.41%	1.537%
40	32.515	5050	5052	5059	rBV7	1358	3082	1.27%	0.145%

Sum of corrected areas: 2122028

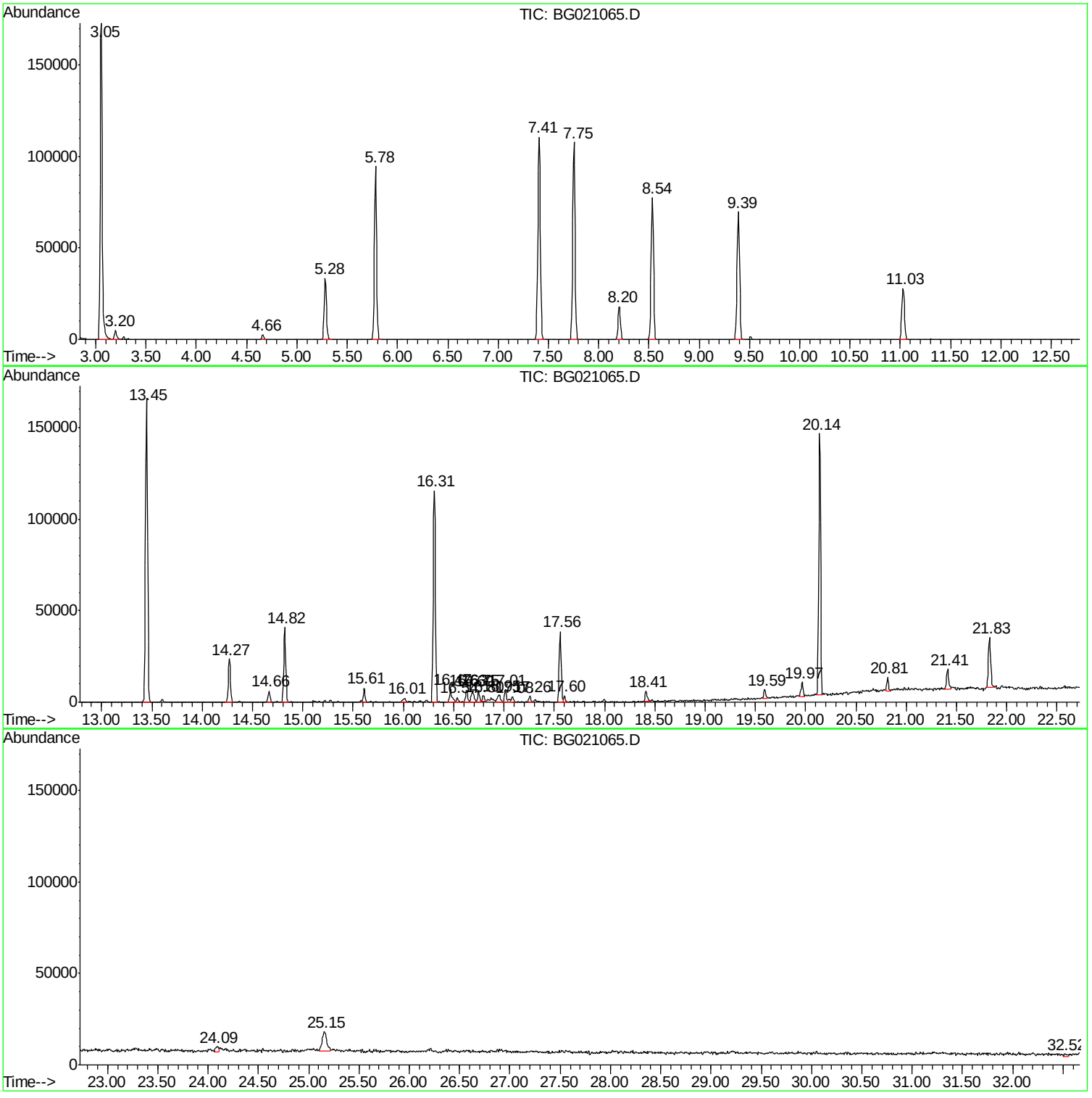
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.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 G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

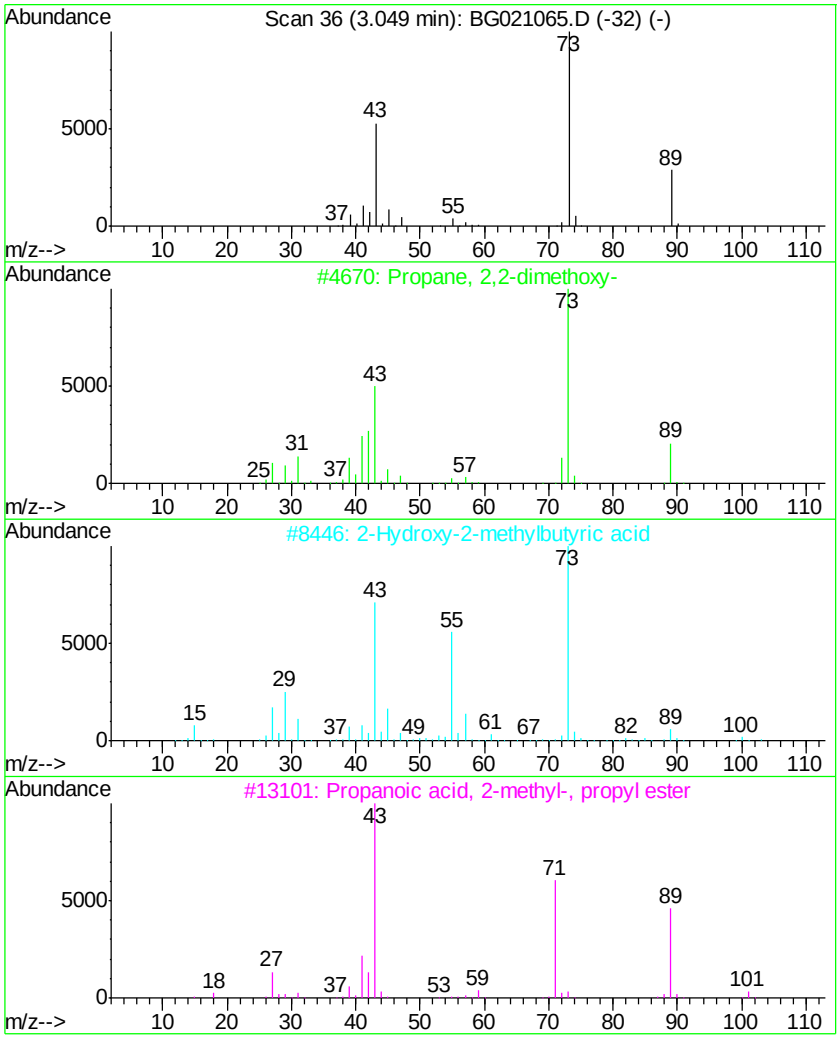
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

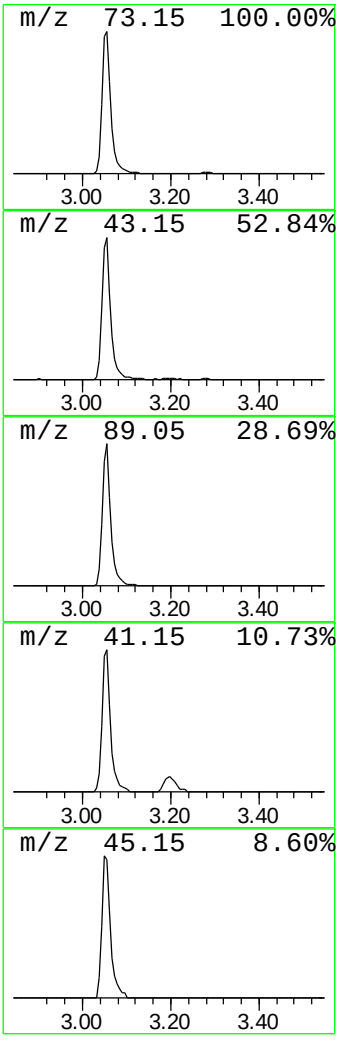
****

Peak Number 1 unknown3.05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.05	167.40 ng	238641	1,4-Dichlorobenzene-d4	8.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	40
2	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
3	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
4	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	7





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

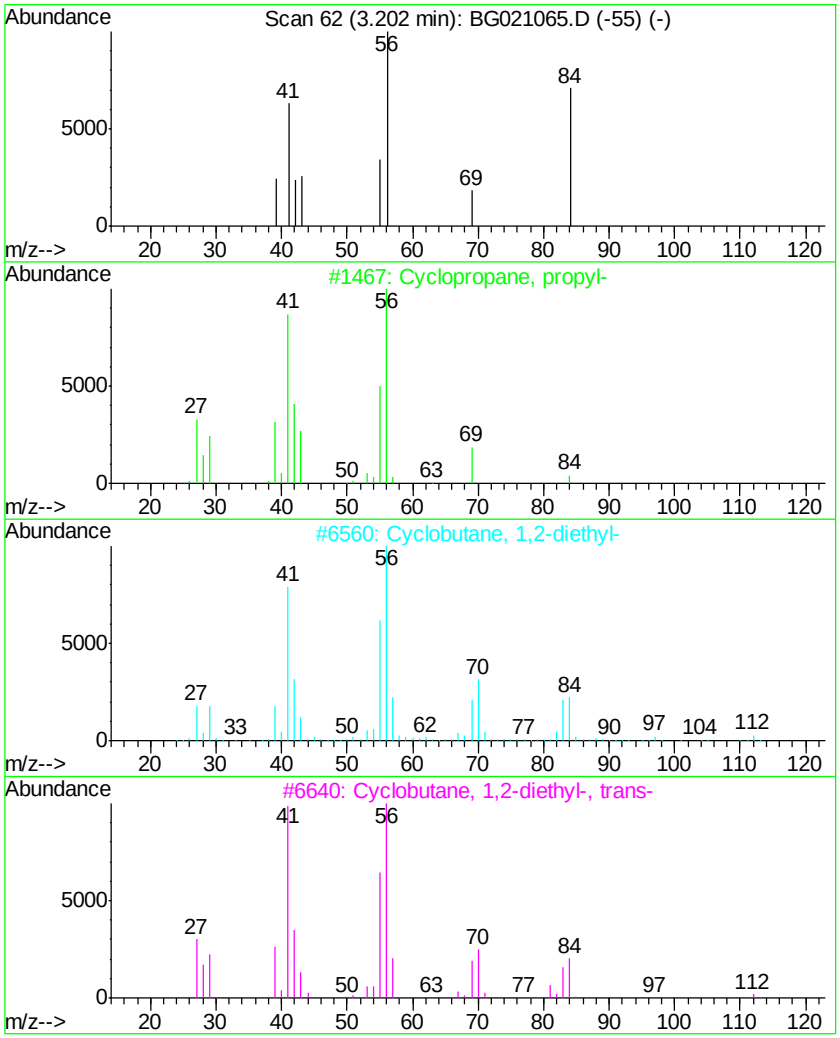
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

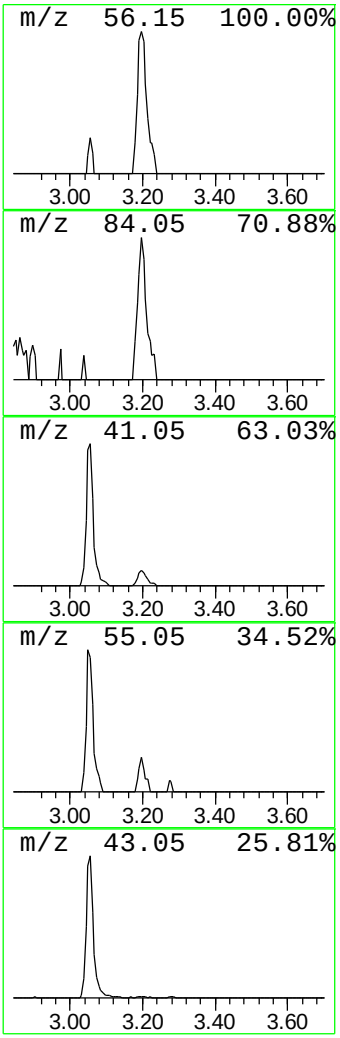
****

Peak Number 2 unknown3.20 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.20	5.90 ng	8408	1,4-Dichlorobenzene-d4	8.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, propyl-	84	C6H12	002415-72-7	38
2	Cyclobutane, 1,2-diethyl-	112	C8H16	061141-83-1	33
3	Cyclobutane, 1,2-diethyl-, trans-	112	C8H16	019341-98-1	9
4	1-Hexanol	102	C6H14O	000111-27-3	9
5	Pentane, 3-methylene-	84	C6H12	000760-21-4	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

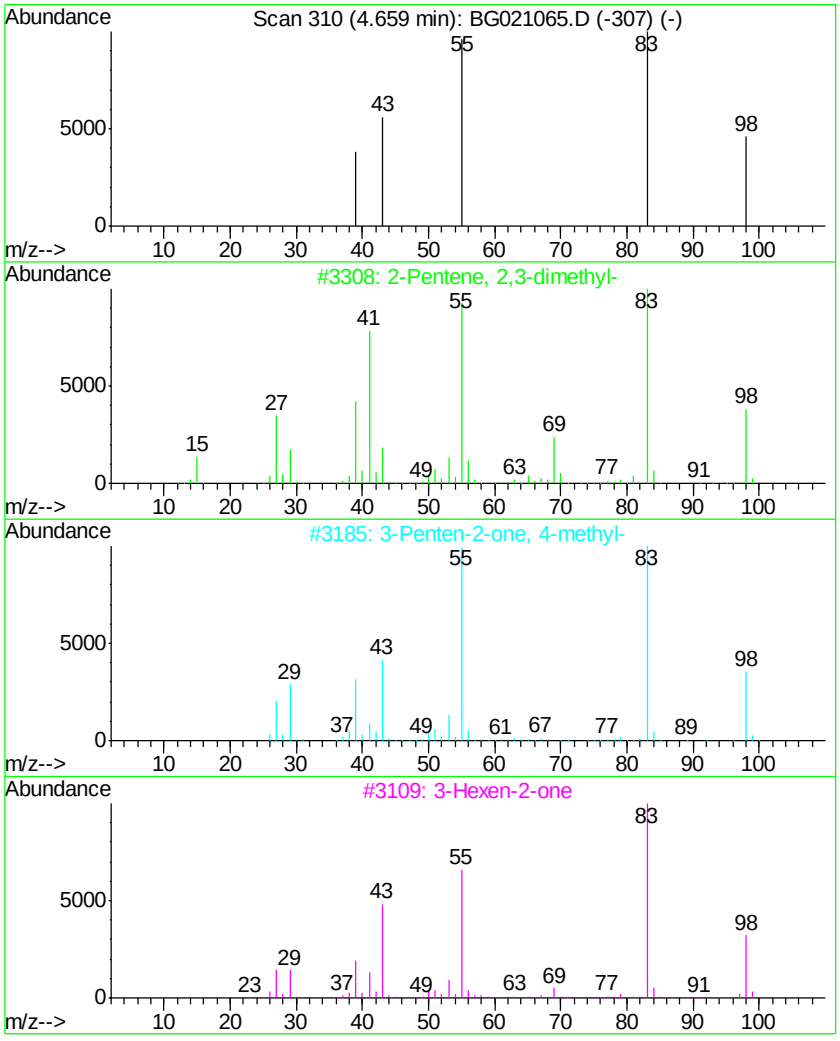
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

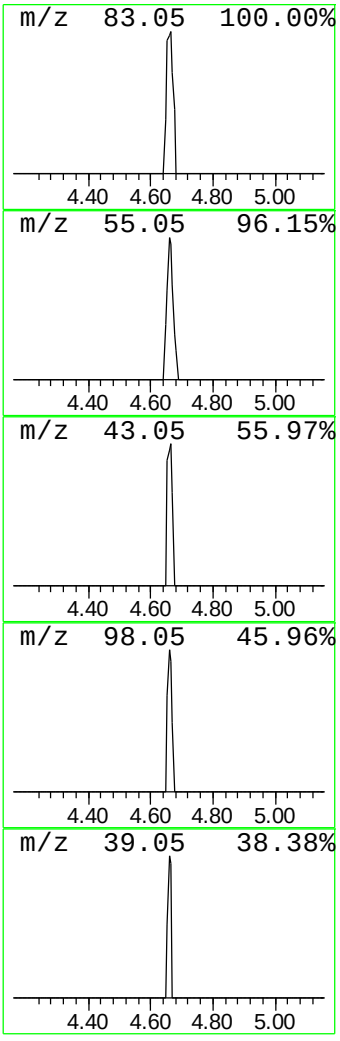
****

Peak Number 3 2-Pentene, 2,3-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	2.50 ng	3564	1,4-Dichlorobenzene-d4	8.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	83
2	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	74
3	3-Hexen-2-one	98	C6H10O	000763-93-9	74
4	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	9
5	2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

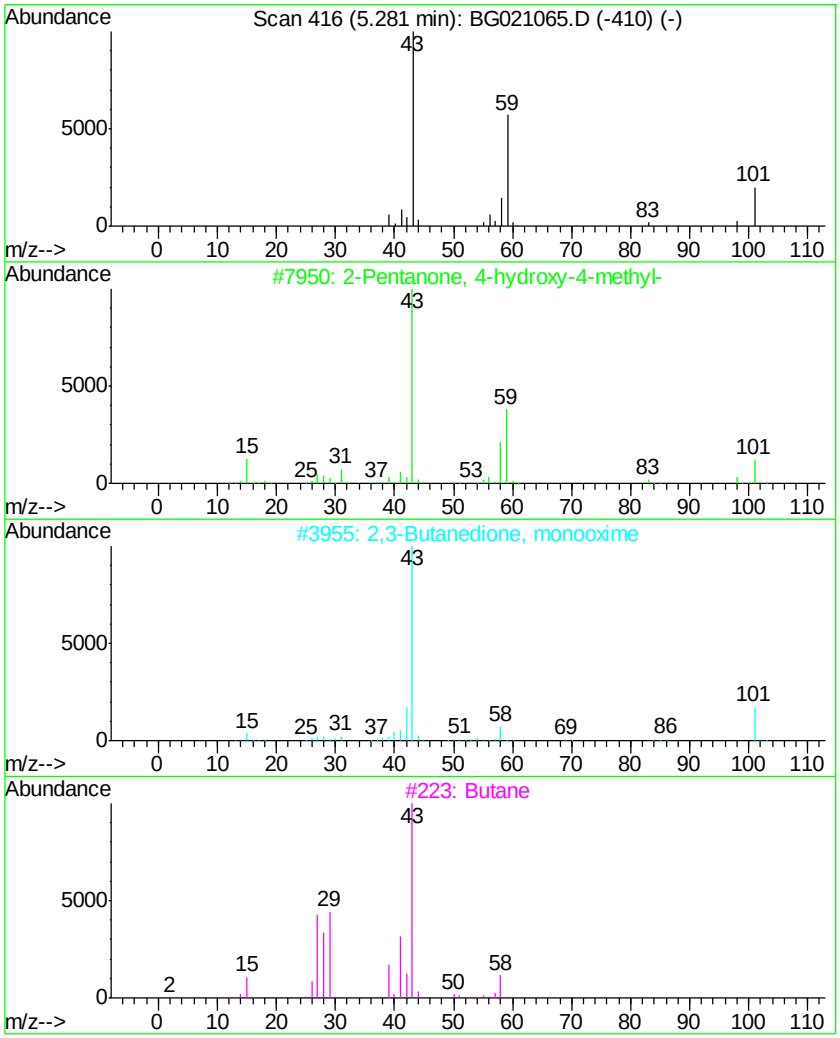
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

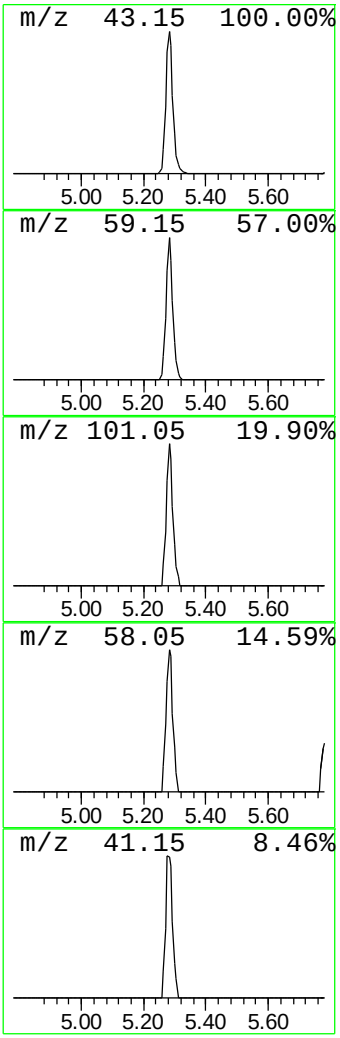
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.28	35.86 ng	51127	1,4-Dichlorobenzene-d4	8.21

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3	Butane	58	C4H10	000106-97-8	9
4	3-Heptanol	116	C7H16O	000589-82-2	9
5	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

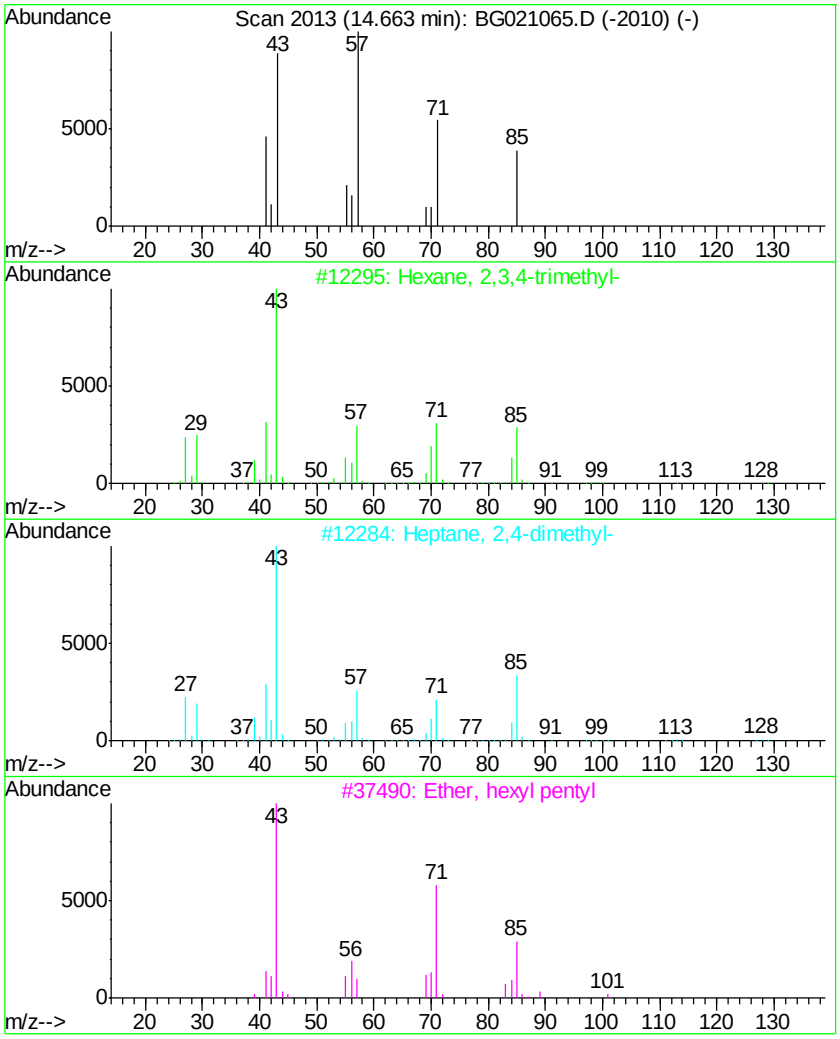
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

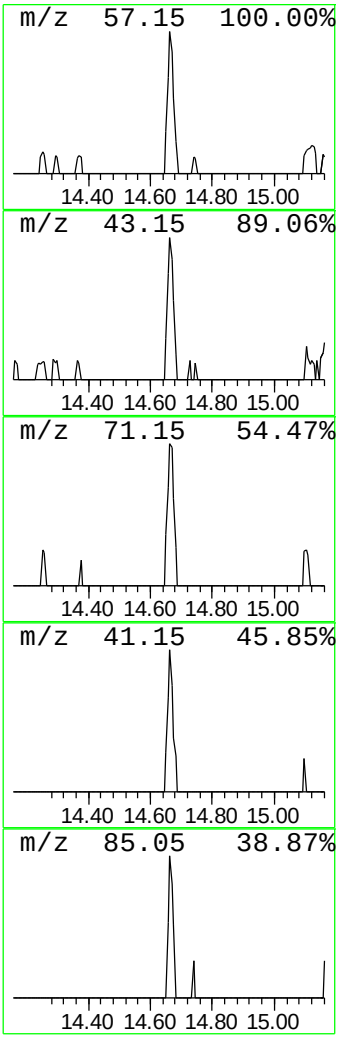
****

Peak Number 7 Hexane, 2,3,4-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	2.47 ng	7508	Acenaphthene-d10	14.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	72
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	42
3	Ether, hexyl pentyl	172	C11H24O	032357-83-8	40
4	Decane, 2,2,6-trimethyl-	184	C13H28	062237-97-2	38
5	Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	38





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

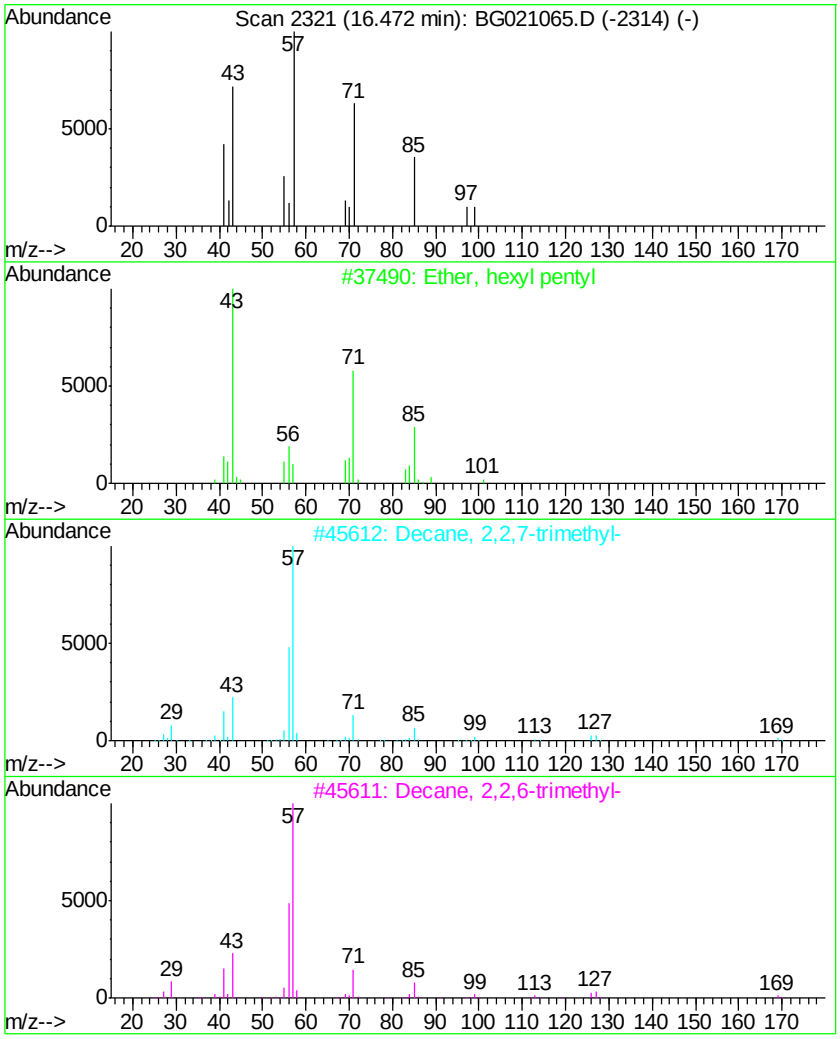
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

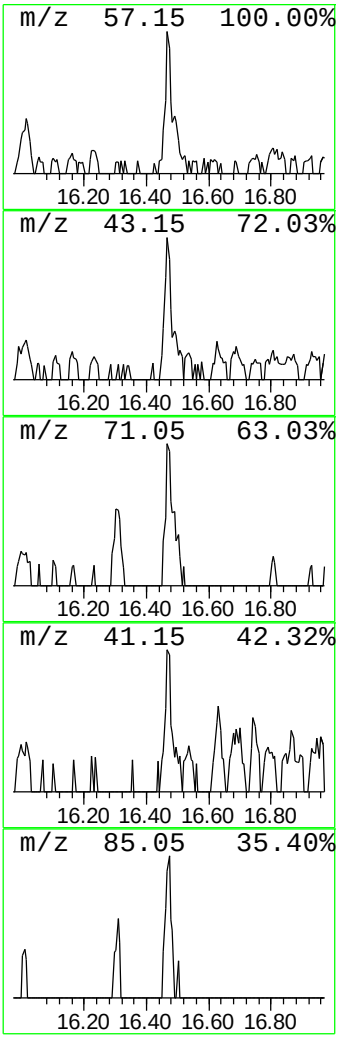
****

Peak Number 9 Ether, hexyl pentyl Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	4.96 ng	13356	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ether, hexyl pentyl	172	C11H24O	032357-83-8	56
2	Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	38
3	Decane, 2,2,6-trimethyl-	184	C13H28	062237-97-2	38
4	2,2-Dimethylpropionic acid, 4-me...	186	C11H22O2	150588-62-8	38
5	Ethanedioic acid, bis(3-methylbu...	230	C12H22O4	002051-00-5	33





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

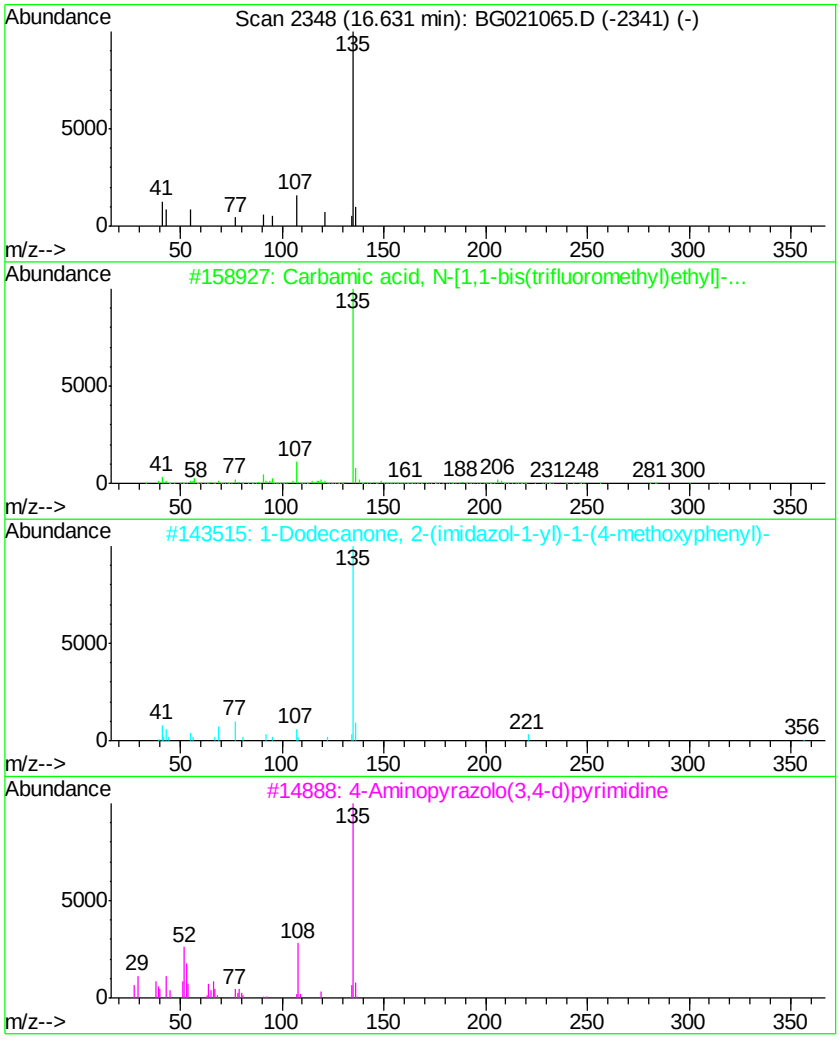
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

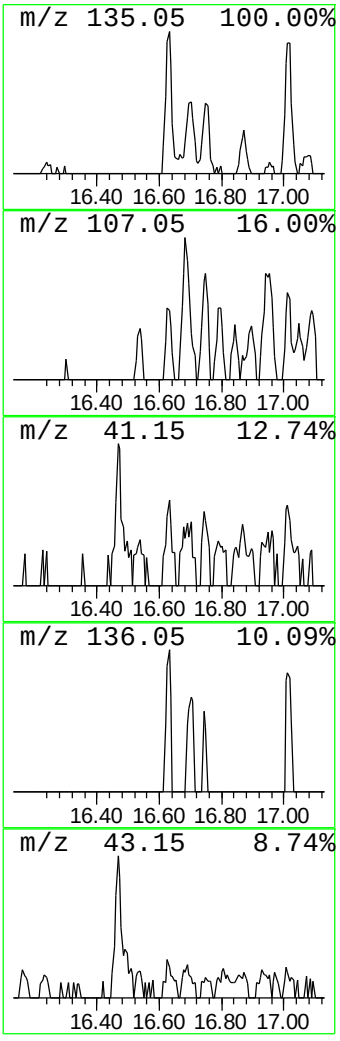
****

Peak Number 10 unknown16.63 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.63	3.63 ng	9771	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6NO2	296242-69-8	43
2	1-Dodecanone, 2-(imidazol-1-yl)-...	356	C22H32N2O2	1000137-95-7	40
3	4-Aminopyrazolo(3,4-d)pyrimidine	135	C5H5N5	020289-44-5	9
4	6-Methyl-triazolo(2,3-b)(1,2,4)-...	135	C5H5N5	061139-75-1	9
5	2H-Imidazo[4,5-b]pyridin-2-one, ...	135	C6H5N3O	016328-62-4	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

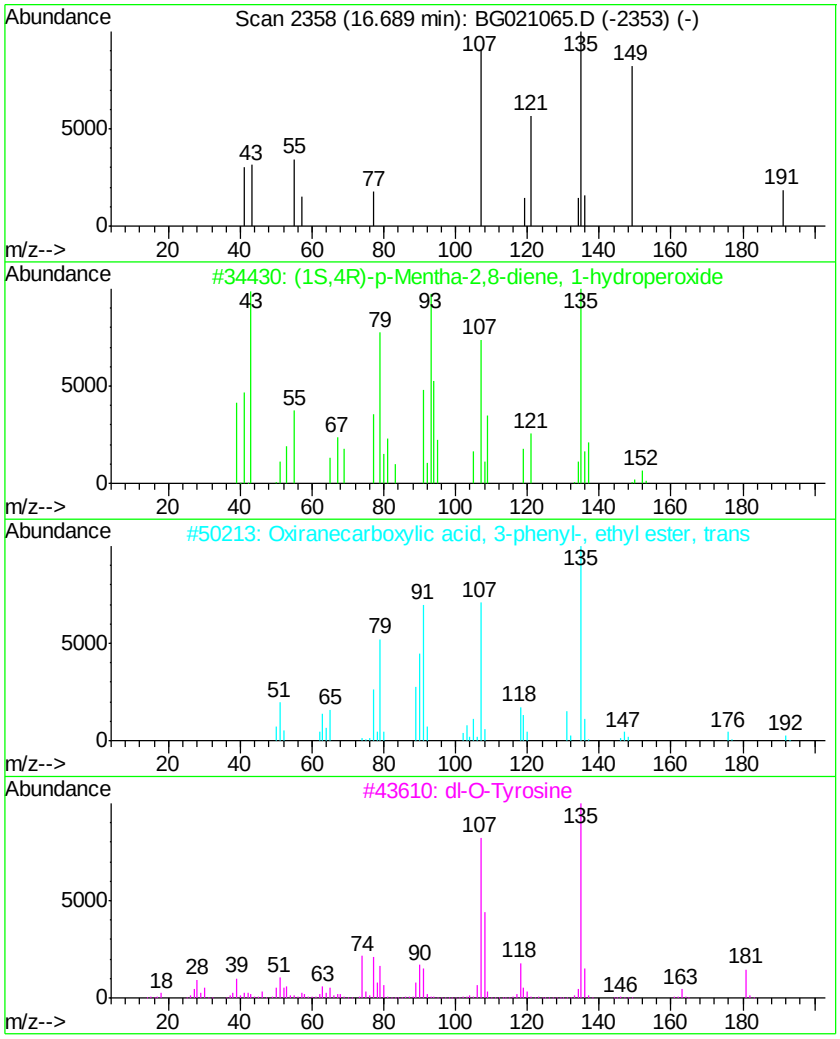
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

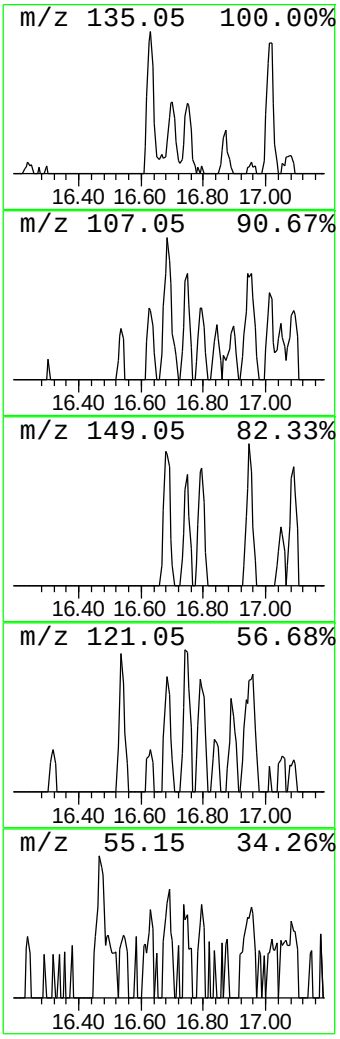
****

Peak Number 11 unknown16.69 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	3.83 ng	10311	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1S,4R)-p-Mentha-2,8-diene, 1-hy...	168	C10H16O2	1000292-73-9	33
2	Oxiranecarboxylic acid, 3-phenyl...	192	C11H12O3	002272-55-1	28
3	dl-O-Tyrosine	181	C9H11NO3	002370-61-8	28
4	Pyridine, 4-ethyl-2,6-dimethyl-	135	C9H13N	036917-36-9	16
5	Benzamide, 3-methoxy-N-allyl-	191	C11H13NO2	331989-39-0	16





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.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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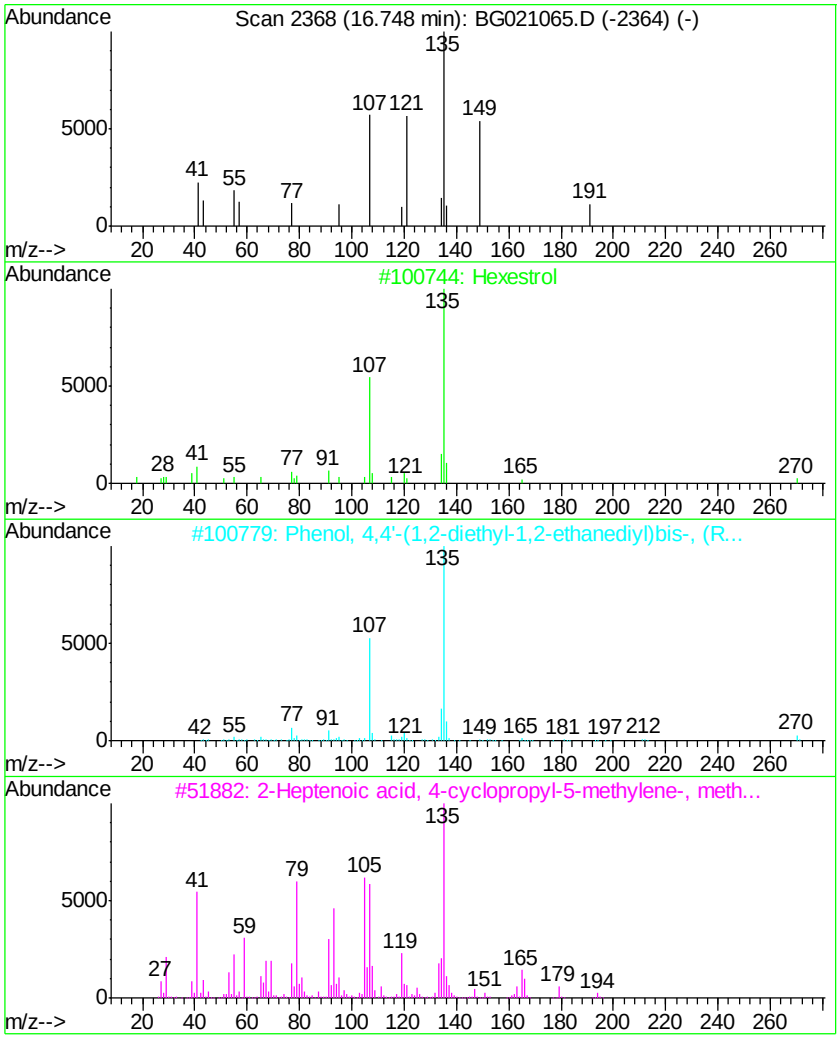
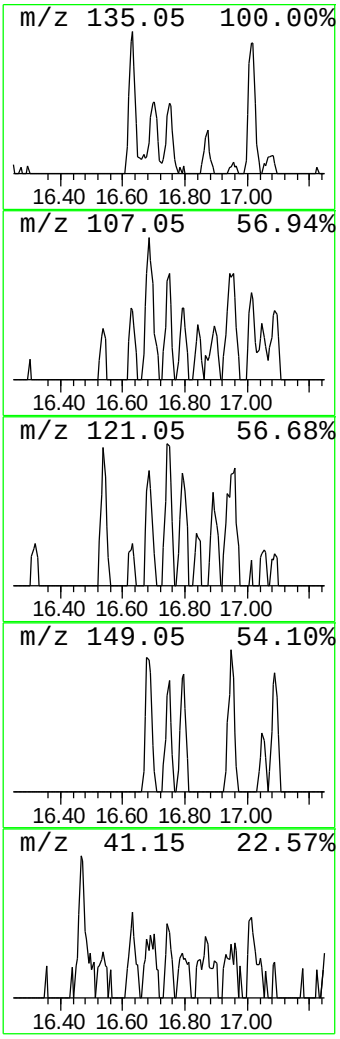
Peak Number 12 Hexestrol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.75	3.06 ng	8234	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexestrol	270	C18H22O2	005635-50-7	50
2	Phenol, 4,4'-(1,2-diethyl-1,2-et...	270	C18H22O2	000084-16-2	40
3	2-Heptenoic acid, 4-cyclopropyl-...	194	C12H18O2	074793-23-0	33
4	1H-Indole, 5-fluoro-	135	C8H6FN	000399-52-0	32
5	Oxiranecarboxylic acid, 3-phenyl...	192	C11H12O3	002272-55-1	28





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022416\
Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.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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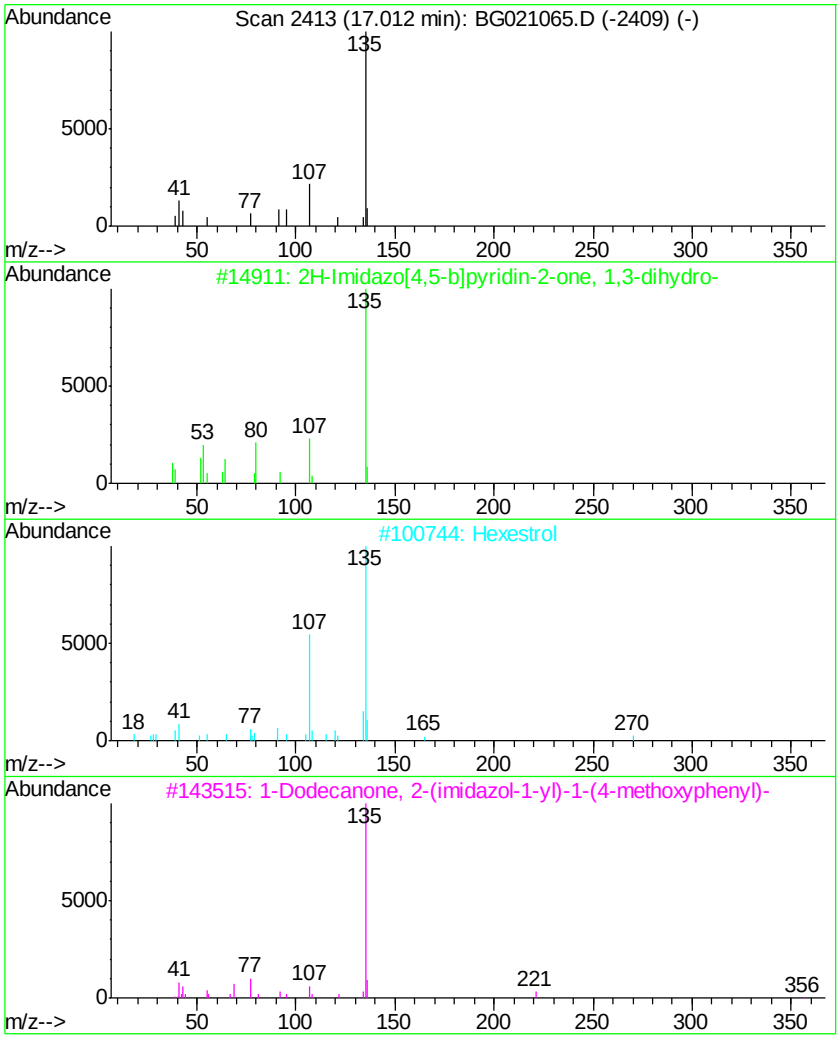
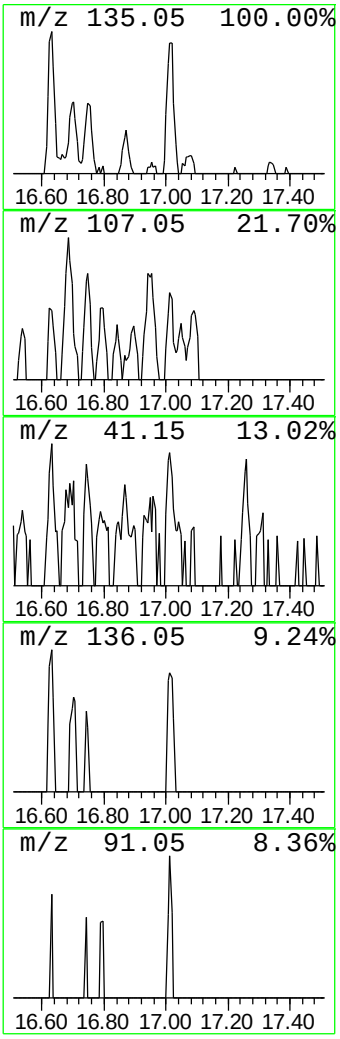
Peak Number 14 2H-Imidazo[4,5-b]pyridin-2-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.01	3.50 ng	9430	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Imidazo[4,5-b]pyridin-2-one, ...	135	C6H5N3O	016328-62-4	50
2	Hexestrol	270	C18H22O2	005635-50-7	39
3	1-Dodecanone, 2-(imidazol-1-yl)-...	356	C22H32N2O2	1000137-95-7	38
4	Benzeneacetonitrile, 2-fluoro-	135	C8H6FN	000326-62-5	9
5	Formamide, N-methyl-N-phenyl-	135	C8H9NO	000093-61-8	9





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Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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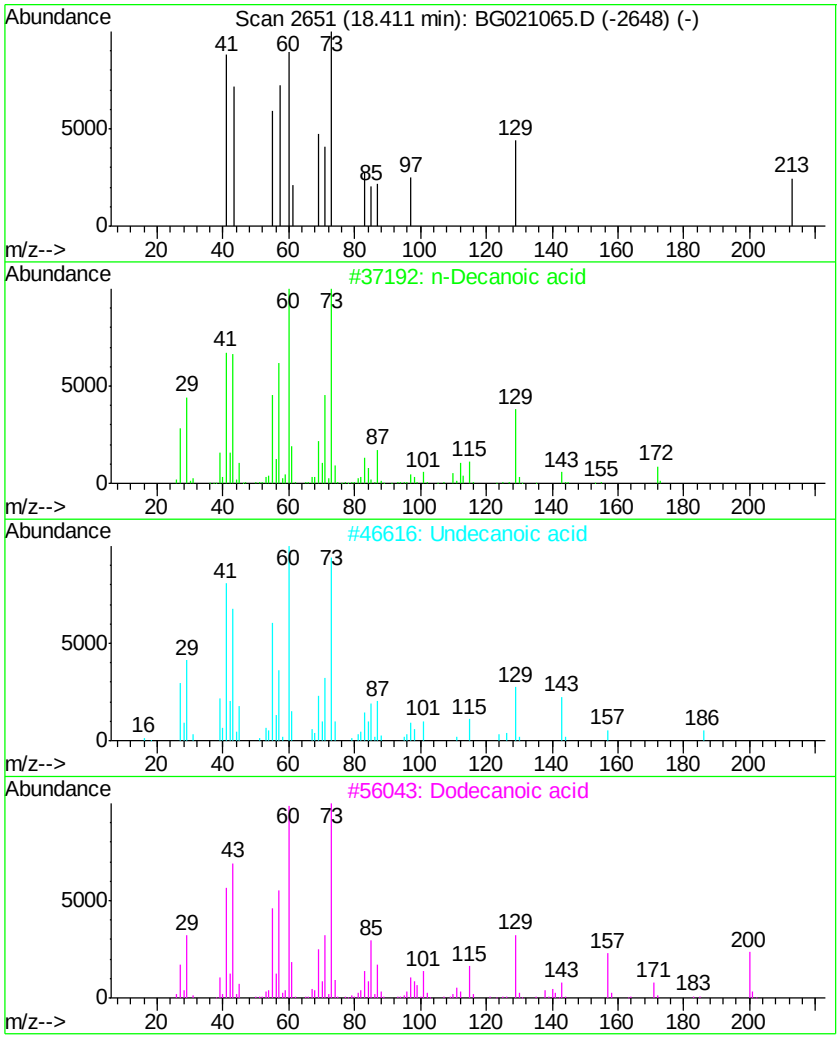
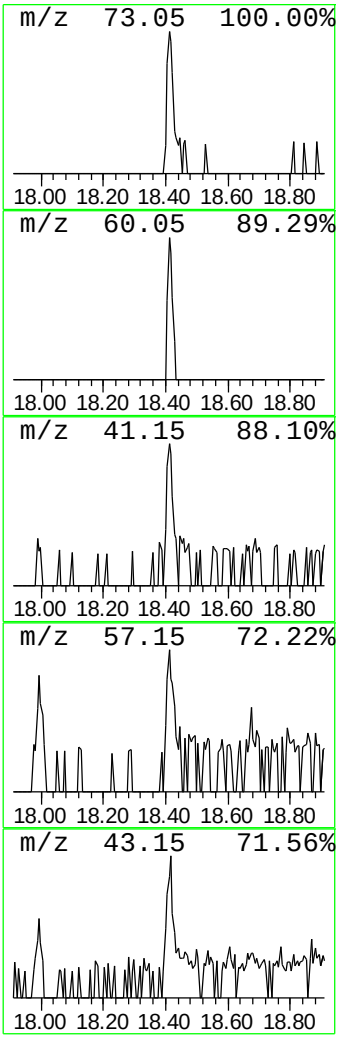
Peak Number 15 n-Decanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	3.11 ng	8376	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	59
2	Undecanoic acid	186	C11H22O2	000112-37-8	53
3	Dodecanoic acid	200	C12H24O2	000143-07-7	45
4	Lactose	342	C12H22O11	000063-42-3	38
5	1-Fluorononane	146	C9H19F	000463-18-3	27





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Data File : BG021065.D
Acq On : 24 Feb 2016 23:51
Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
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TIC Integration Parameters: LSCINT.P

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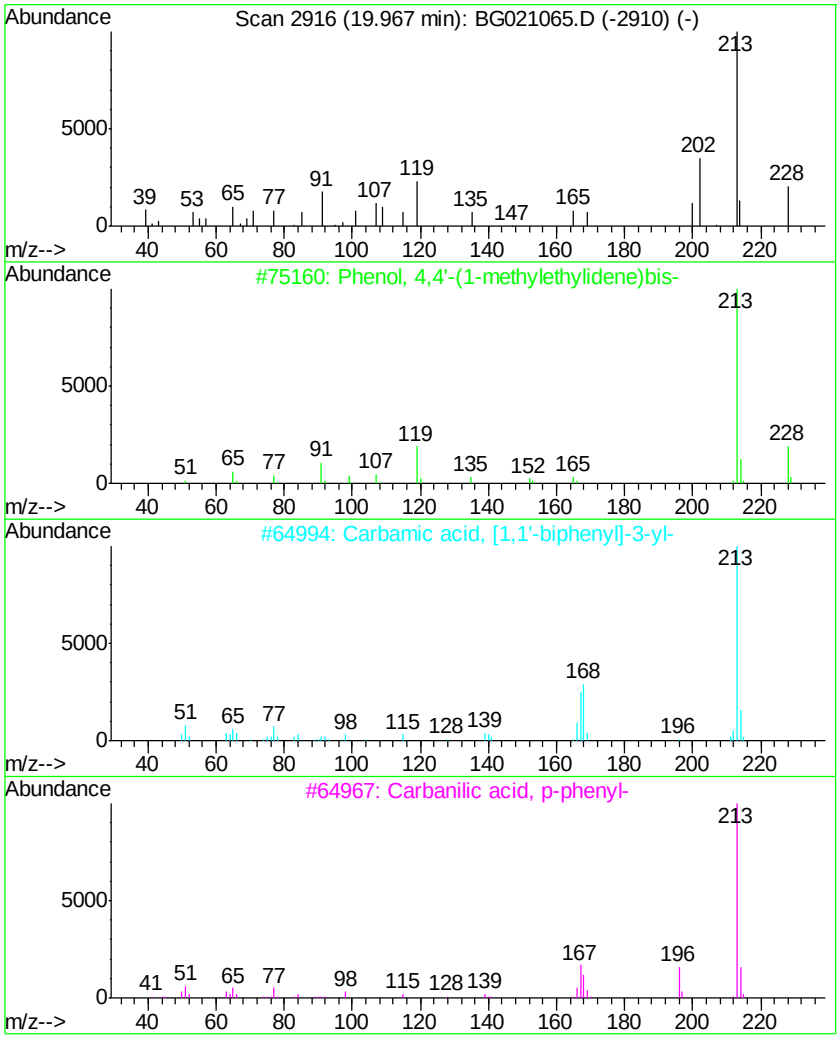
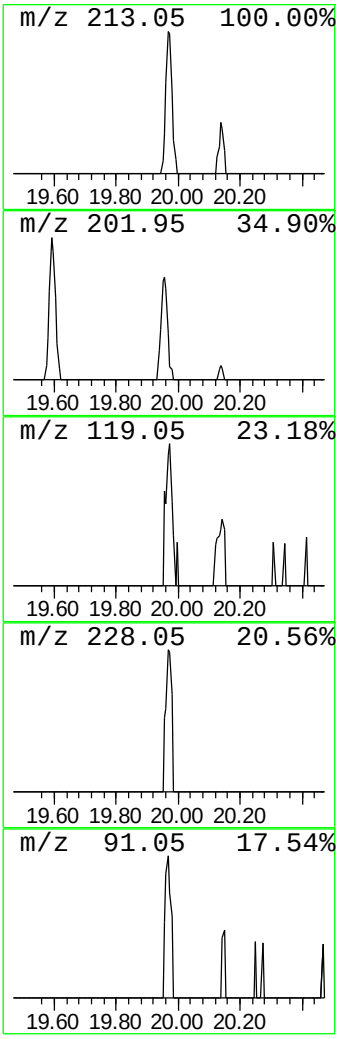
Peak Number 16 Phenol, 4,4'-(1-methylethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.97	5.44 ng	11459	Chrysene-d12	21.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	64
2	Carbamic acid, [1,1'-biphenyl]-3...	213	C13H11NO2	055030-29-0	32
3	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	32
4	2-Trifluoromethyl-5-hydroxyquino...	213	C10H6F3NO	041958-06-9	32
5	3-Dibenzofuranamine, 2-methoxy-	213	C13H11NO2	005834-17-3	25





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Operator : UM/SJ
Sample : H1521-11
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SP-3+20(0-2)

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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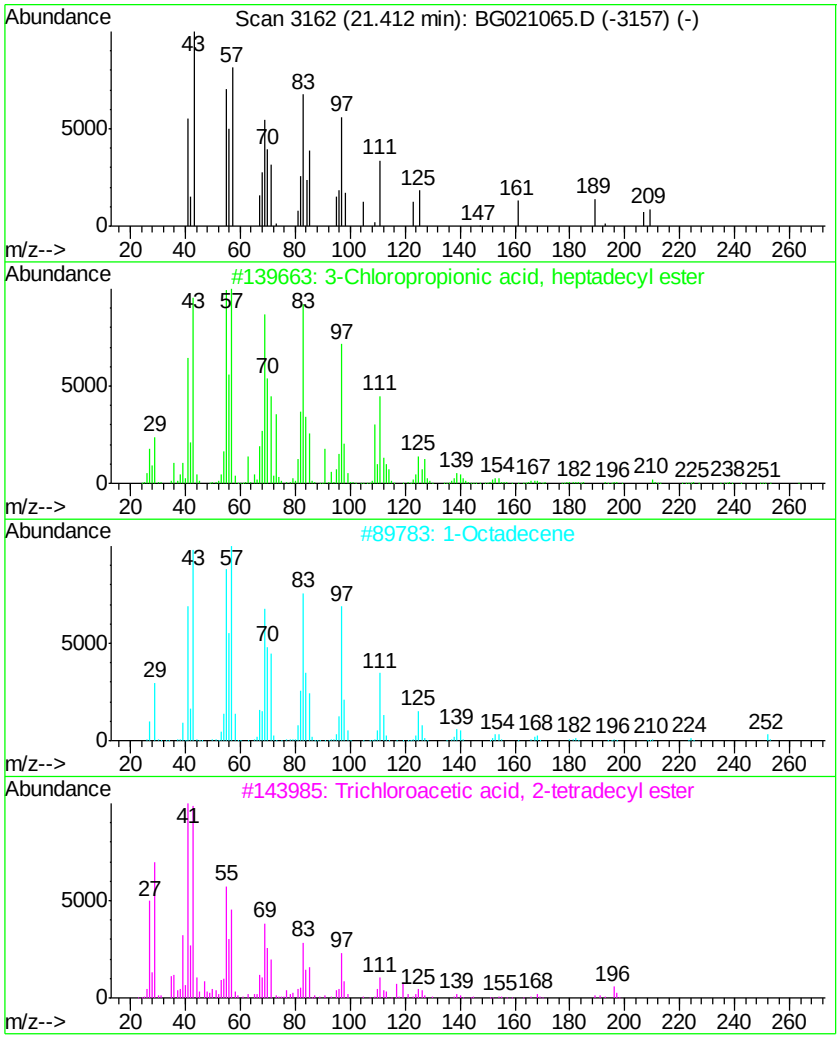
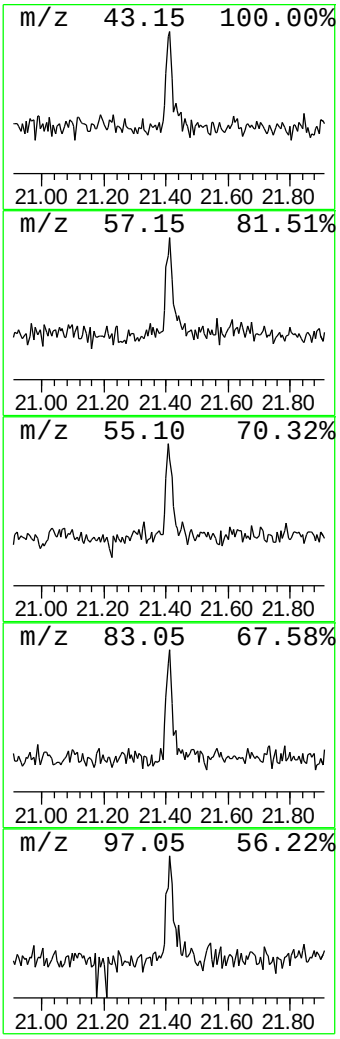
Peak Number 18 3-Chloropropionic acid, hep... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.41	7.83 ng	16492	Chrysene-d12	21.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	74
2	1-Octadecene	252	C18H36	000112-88-9	74
3	Trichloroacetic acid, 2-tetradec...	358	C16H29Cl3O2	1000282-06-6	72
4	Hexadecen-1-ol, trans-9-	240	C16H32O	064437-47-4	72
5	9-Eicosene, (E)-	280	C20H40	074685-29-3	72





Tentatively Identified Compound (LSC) summary

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Data File : BG021065.D
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Operator : UM/SJ
Sample : H1521-11
Misc :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown3.05	3.05	167.4	ng	238641	1	8.21	28512	20.0
unknown3.20	3.20	5.9	ng	8408	1	8.21	28512	20.0
2-Pentene, 2,3-di...	4.66	2.5	ng	3564	1	8.21	28512	20.0
2-Pentanone, 4-hy...	5.28	35.9	ng	51127	1	8.21	28512	20.0
Hexane, 2,3,4-tri...	14.66	2.5	ng	7508	3	14.82	60745	20.0
Ether, hexyl pentyl	16.47	5.0	ng	13356	4	17.56	53809	20.0
unknown16.63	16.63	3.6	ng	9771	4	17.56	53809	20.0
unknown16.69	16.69	3.8	ng	10311	4	17.56	53809	20.0
Hexestrol	16.75	3.1	ng	8234	4	17.56	53809	20.0
2H-Imidazo[4,5-b]...	17.01	3.5	ng	9430	4	17.56	53809	20.0
n-Decanoic acid	18.41	3.1	ng	8376	4	17.56	53809	20.0
Phenol, 4,4'-(1-m...	19.97	5.4	ng	11459	5	21.83	42150	20.0
3-Chloropropionic...	21.41	7.8	ng	16492	5	21.83	42150	20.0