LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

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-BG061215.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	5.115	407	413	424	rVB	211420	325180	10.67%	2.853%
2	6.731	682	688	698	rBV	146978	234617	7.70%	2.058%
3	7.066	737	745	752	rBV	418694	635486	20.85%	5.575%
4	7.524	817	823	828	rBV	95415	143635	4.71%	1.260%
5	7.841	870	877	882	rBV	363314	561000	18.40%	4.922%
6	7.894	882	886	894	rVB2	47659	77068	2.53%	0.676%
7	8.693	1015	1022	1030	rVB	296892	482040	15.81%	4.229%
8	8.963	1063	1068	1072	rBV3	23223	37125	1.22%	0.326%
9	9.527	1159	1164	1169	rBV2	48297	81895	2.69%	0.718%
10	9.721	1188	1197	1204	rBV3	41988	80525	2.64%	0.706%
11	9.845	1209	1218	1228	rVB4	37494	82441	2.70%	0.723%
12	10.321	1293	1299	1302	rBV	105567	171272	5.62%	1.503%
13	10.368	1302	1307	1315	rVB	190077	312589	10.26%	2.742%
14	11.425	1482	1487	1493	rBV4	18658	32853	1.08%	0.288%
15	11.889	1561	1566	1575	rVB4	16272	31604	1.04%	0.277%
16	12.218	1615	1622	1627	rVB5	19517	37679	1.24%	0.331%
17	12.818	1718	1724	1730	rBV	877697	1210628	39.72%	10.621%
18	13.529	1838	1845	1849	rBV5	18898	40800	1.34%	0.358%
19	14.204	1955	1960	1966	rBV2	197995	288393	9.46%	2.530%
20	14.269	1966	1971	1979	rVB	122954	179564	5.89%	1.575%
21	15.262	2136	2140	2146	rBV4	20191	32190	1.06%	0.282%
22	15.708	2208	2216	2224	rBV	968184	1345987	44.16%	11.809%
23	15.861	2239	2242	2246	rBV3	28304	33863	1.11%	0.297%
24	15.938	2251	2255	2262	rVB9	27532	55282	1.81%	0.485%
25	16.960	2423	2429	2434	rBV2	317575	429044	14.08%	3.764%
26	17.871	2580	2584	2588	rVV7	30271	48974	1.61%	0.430%
27	17.947	2592	2597	2603	rVV4	46191	72537	2.38%	0.636%
28	18.793	2736	2741	2746	rVB4	25319	37716	1.24%	0.331%
29	19.604	2873	2879	2889	rBV	2471601	3048136	100.00%	26.742%
30	19.815	2912	2915	2920	rVB5	28737	30890	1.01%	0.271%
31	20.973	3110	3112	3117	rVB6	22469	32283	1.06%	0.283%
32	21.161	3139	3144	3151	rVB	505377	609179	19.99%	5.344%
33	23.358	3511	3518	3528	rVB2	309222	575965	18.90%	5.053%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

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-BG061215.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Sum of corrected areas: 11398440

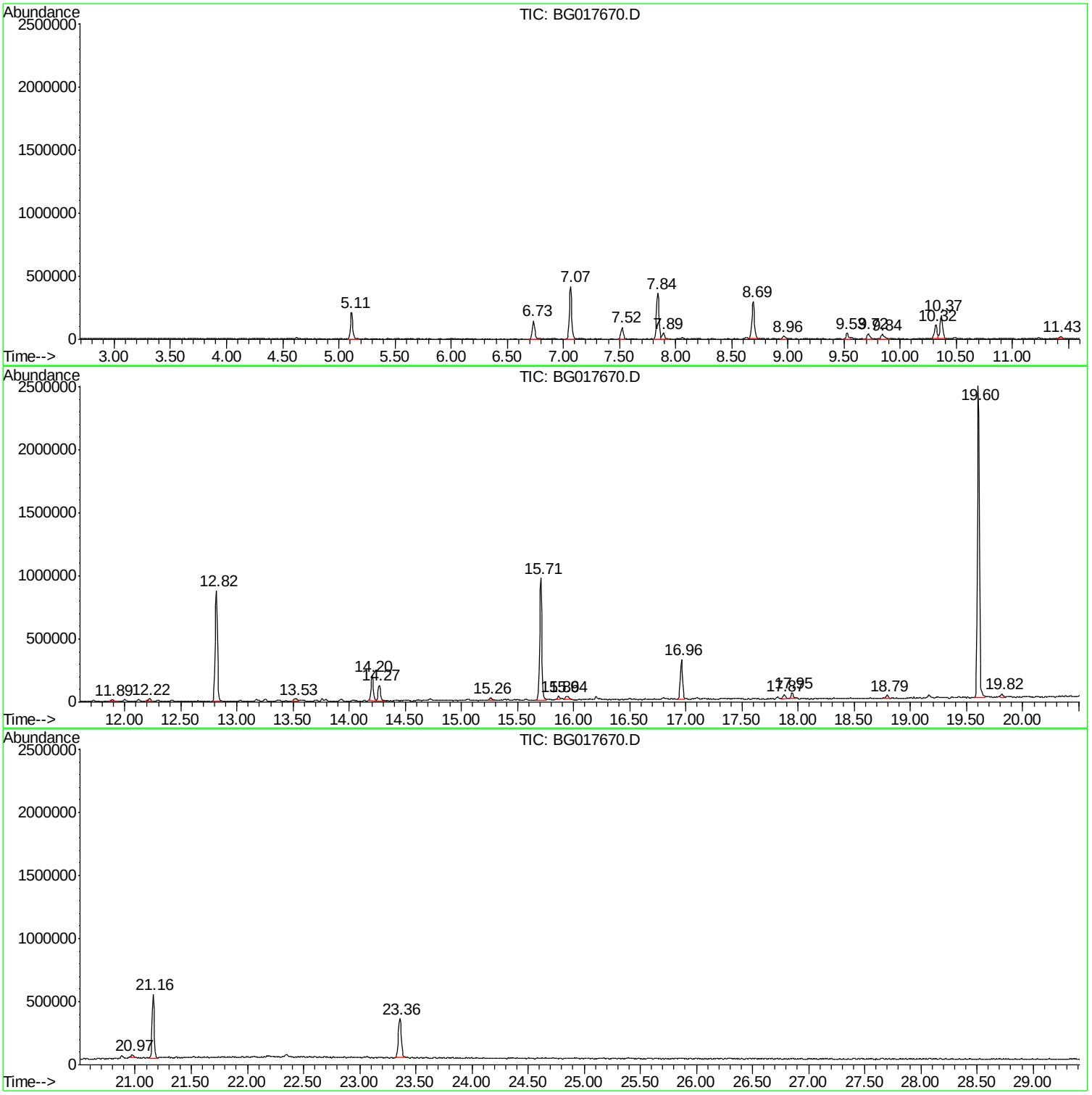
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.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\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

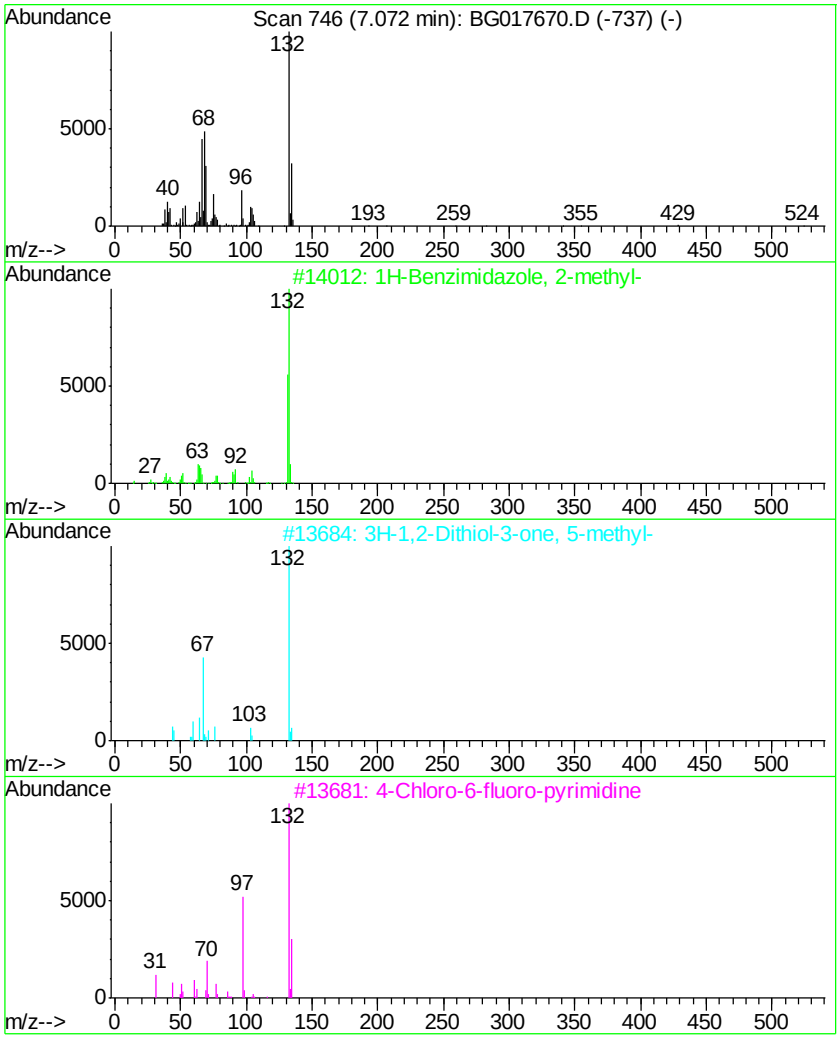
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

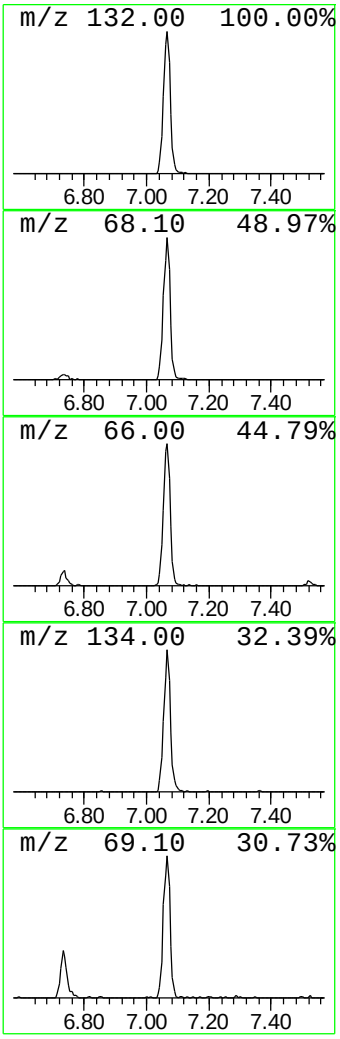
****

Peak Number 1 unknown7.07 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	88.49 ng	635486	1,4-Dichlorobenzene-d4	7.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2	3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
3	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
4	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	12





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

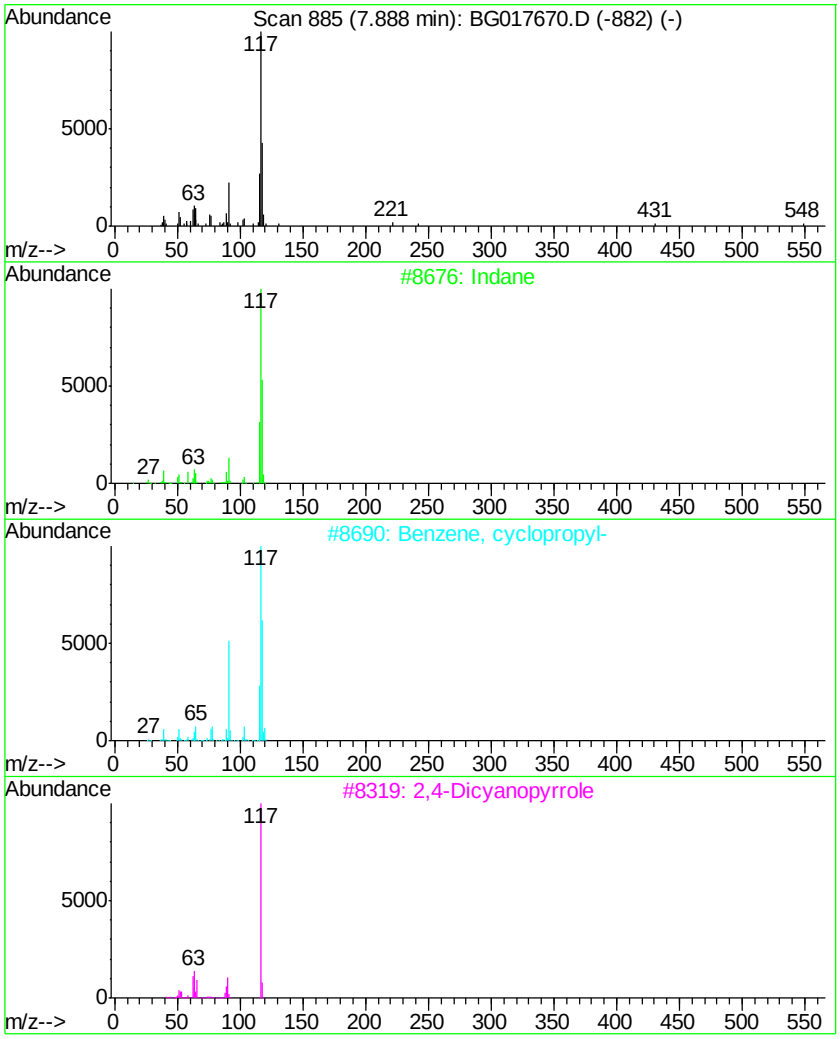
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

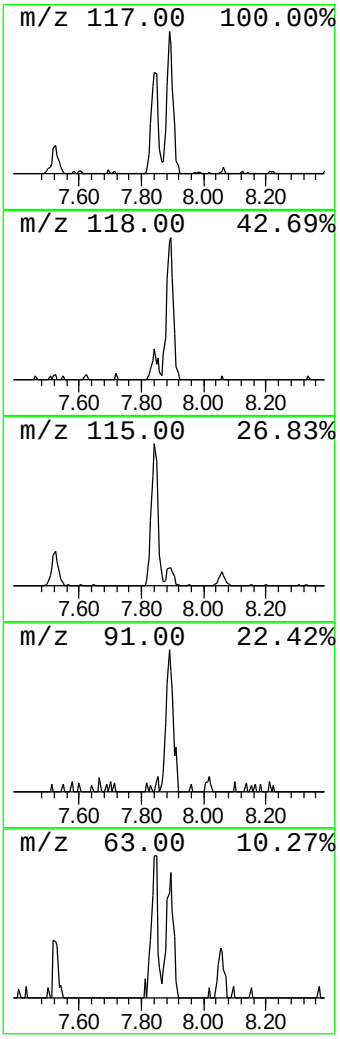
****

Peak Number 3 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.89	10.73 ng	77068	1,4-Dichlorobenzene-d4	7.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	64
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	52
3	2,4-Dicyanopyrrole	117	C6H3N3	074023-87-3	49
4	Deltacyclene	118	C9H10	007785-10-6	49
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	49





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

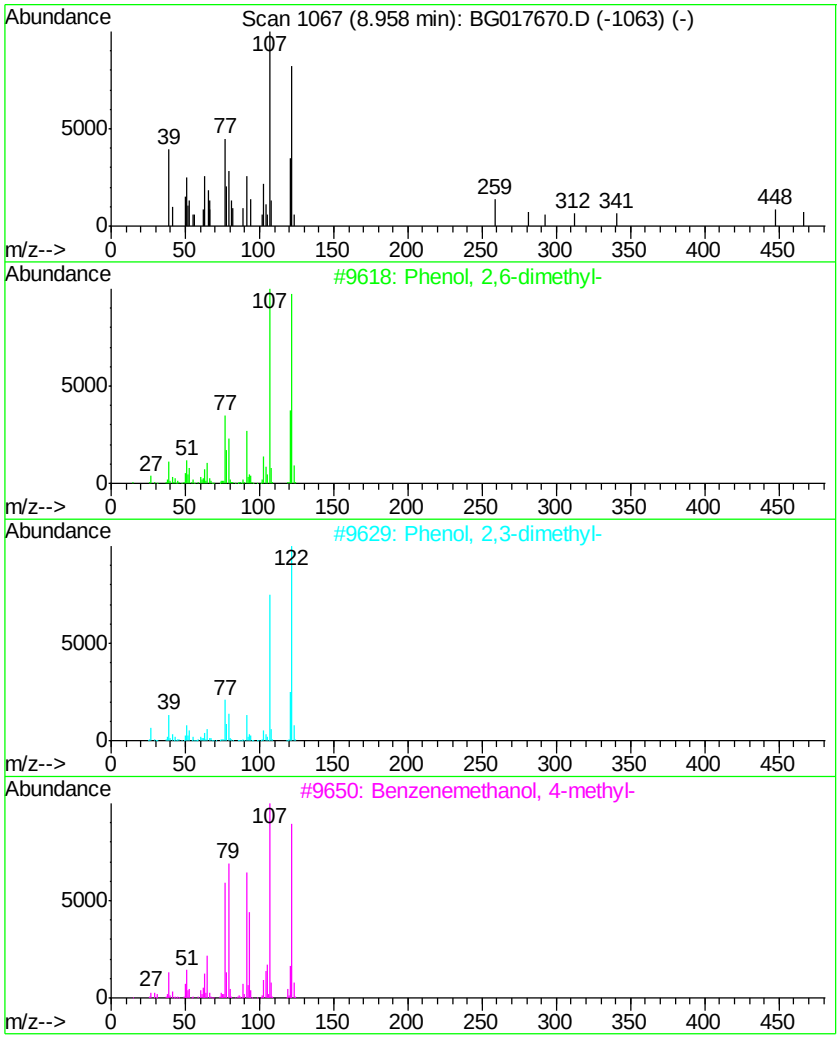
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

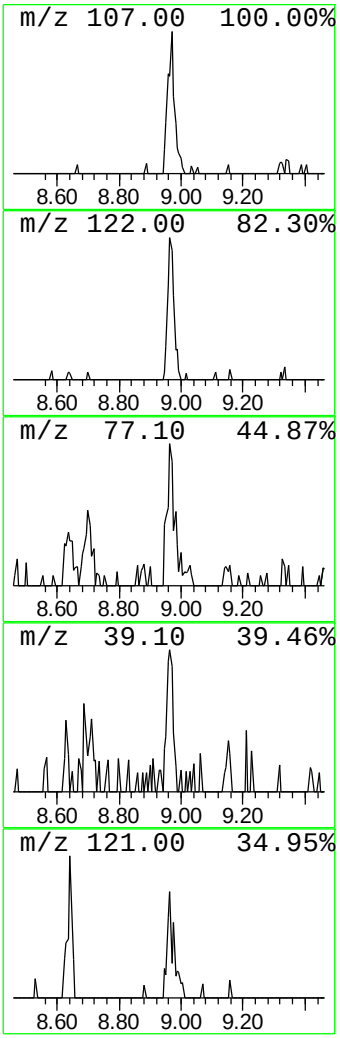
****

Peak Number 4 Phenol, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.96	4.34 ng	37125	Naphthalene-d8	10.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	91
2	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	87
3	Benzenemethanol, 4-methyl-	122	C8H10O	000589-18-4	86
4	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	83
5	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	72





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

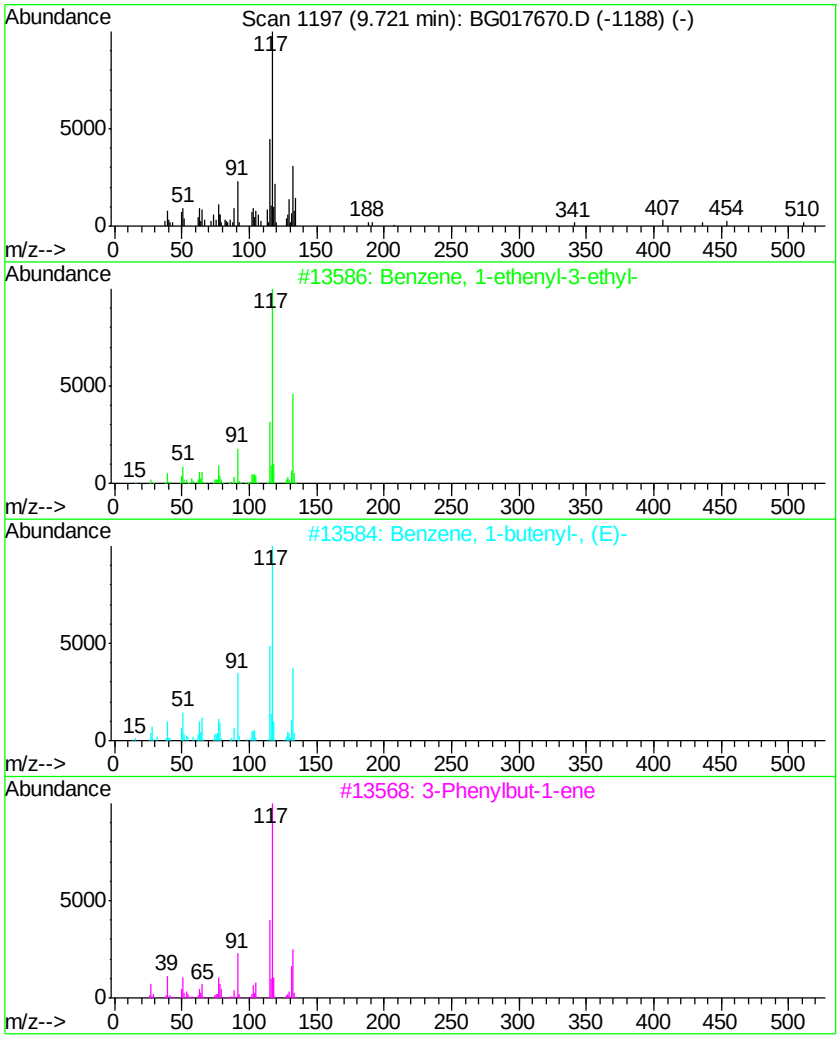
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

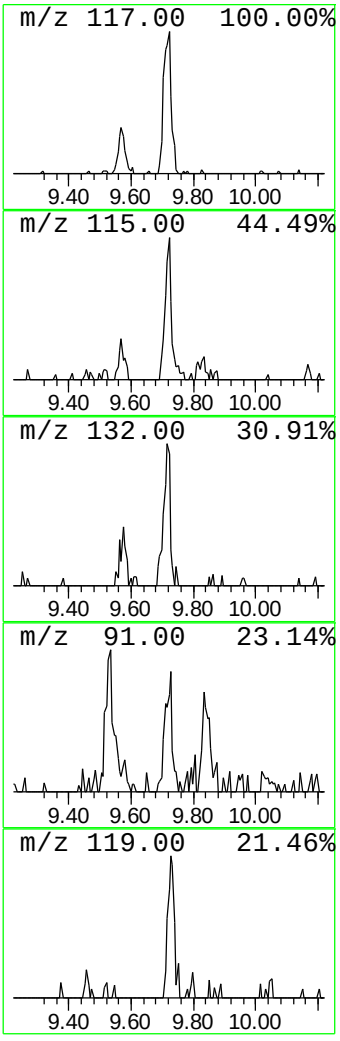
****

Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	9.40 ng	80525	Naphthalene-d8	10.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	89
2	Benzene, 1-butenyl-, (E)-	132	C10H12	001005-64-7	76
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	68
5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	64





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

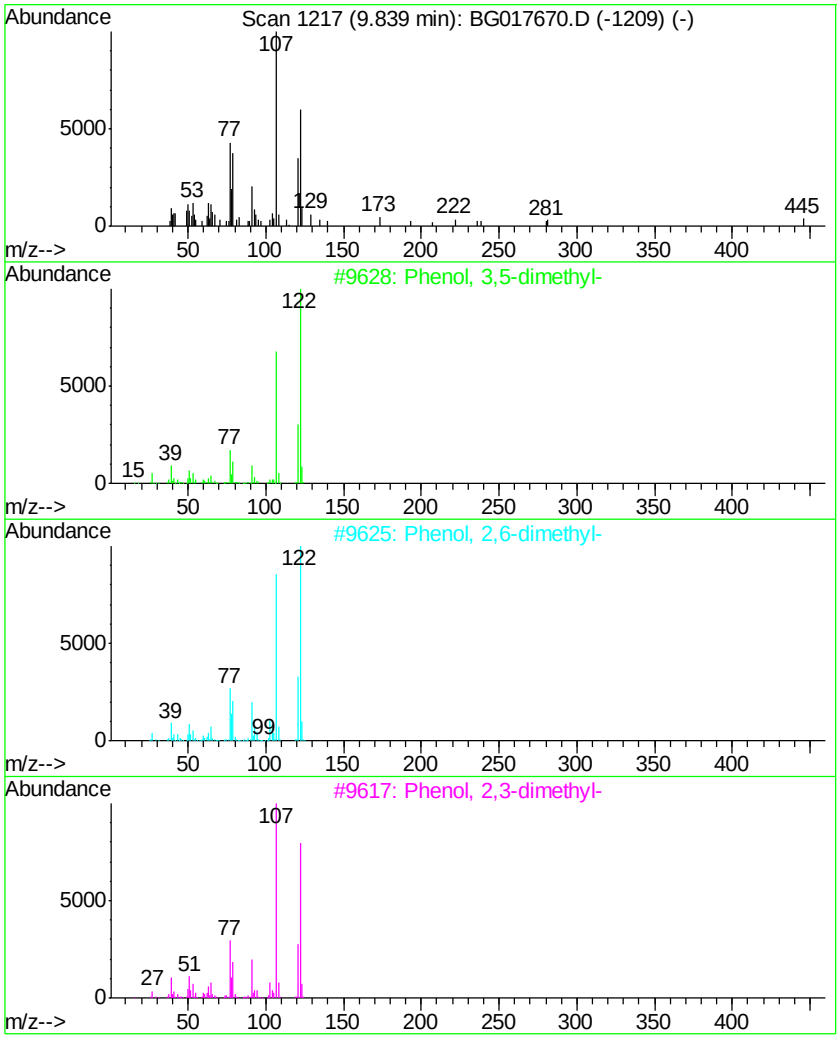
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

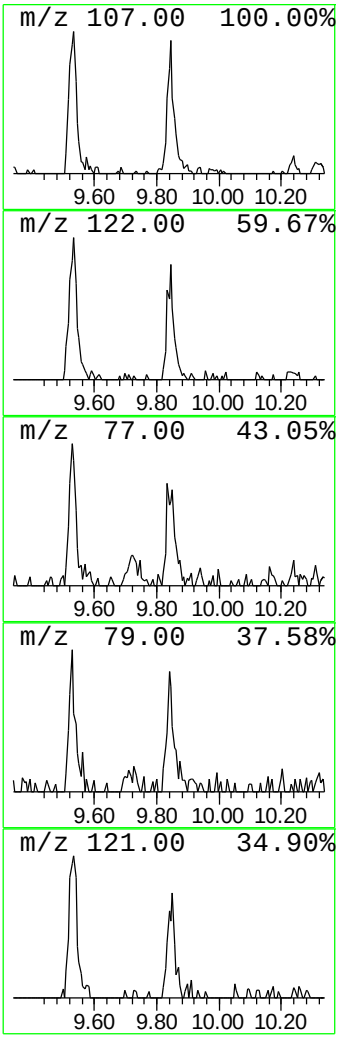
****

Peak Number 6 Phenol, 3,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.84	9.63 ng	82441	Naphthalene-d8	10.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	90
2	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	87
3	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	87
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	80
5	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	80





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

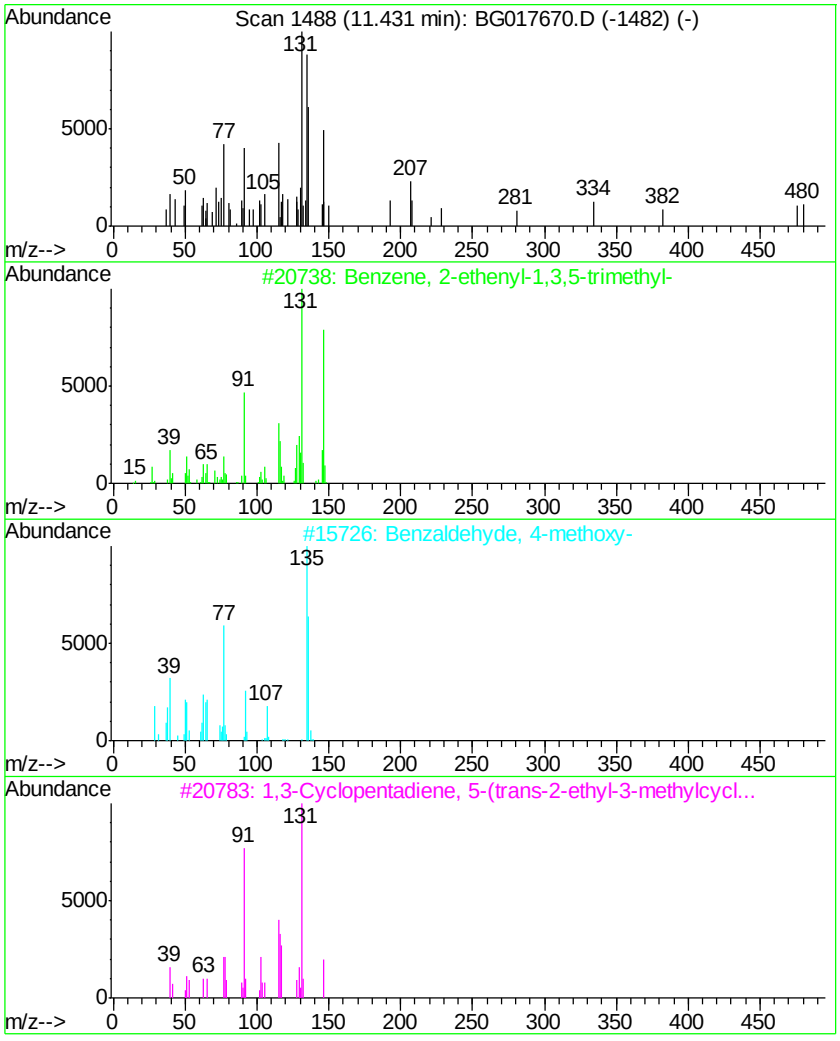
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

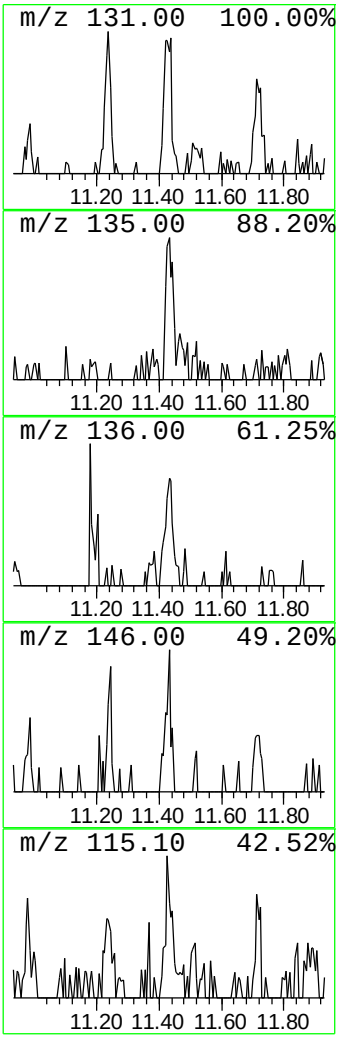
****

Peak Number 7 unknown11.43 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	3.84 ng	32853	Naphthalene-d8	10.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	45
2	Benzaldehyde, 4-methoxy-	136	C8H8O2	000123-11-5	38
3	1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	30
4	4-Methoxymethyl-4-phenyl-1-butene	176	C12H16O	1000215-93-6	27
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	25





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

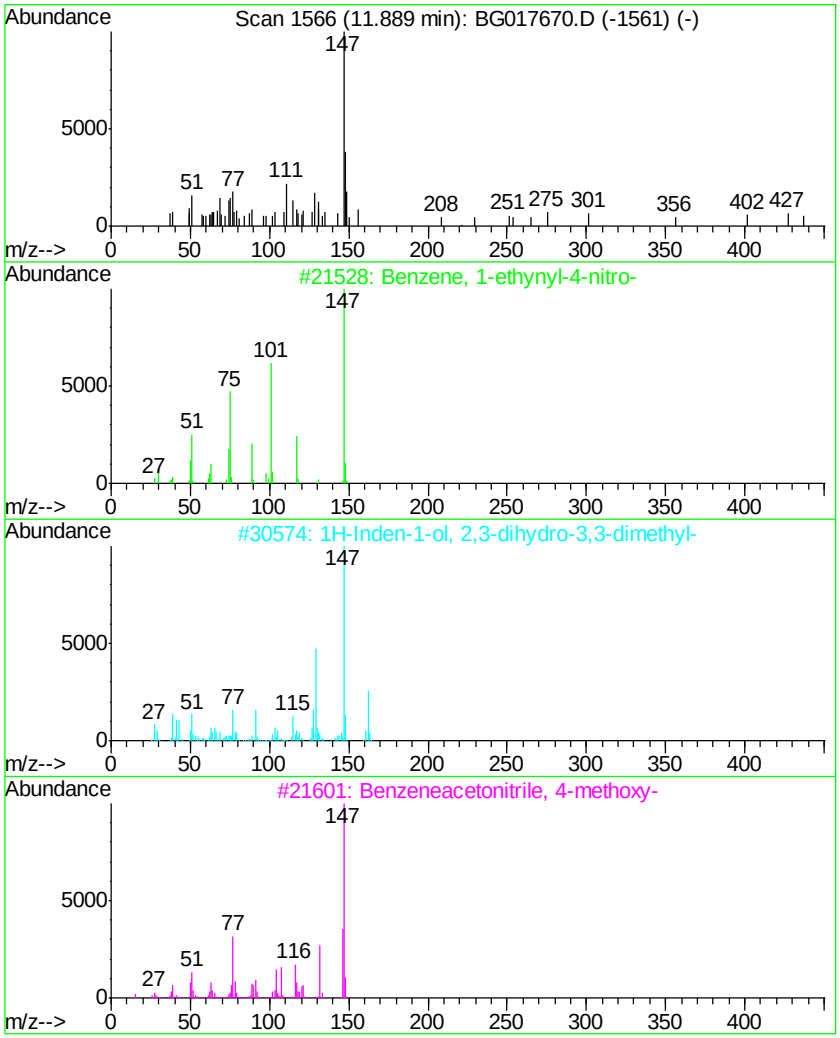
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

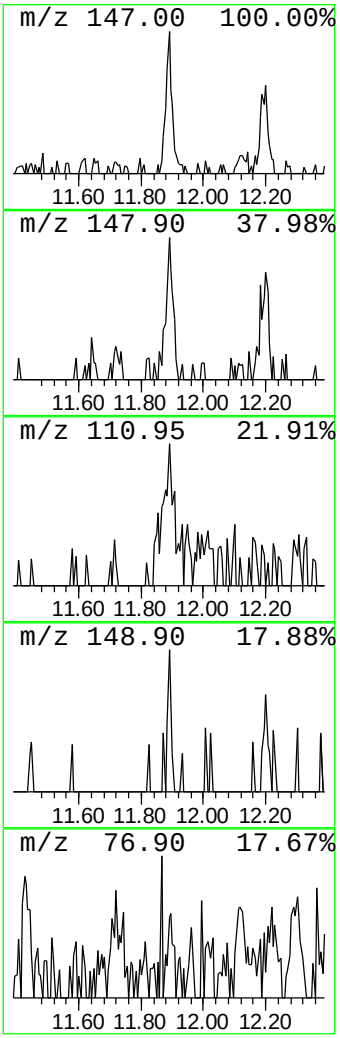
****

Peak Number 8 unknown11.89 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	3.69 ng	31604	Naphthalene-d8	10.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethynyl-4-nitro-	147	C8H5NO2	000937-31-5	46
2	1H-Inden-1-ol, 2,3-dihydro-3,3-d...	162	C11H14O	038393-92-9	43
3	Benzeneacetonitrile, 4-methoxy-	147	C9H9NO	000104-47-2	43
4	Benzene, 1,2-bis(1-methylethyl)-	162	C12H18	000577-55-9	43
5	Ethanone, 1-(2,4,6-trimethylphen...	162	C11H14O	001667-01-2	43





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

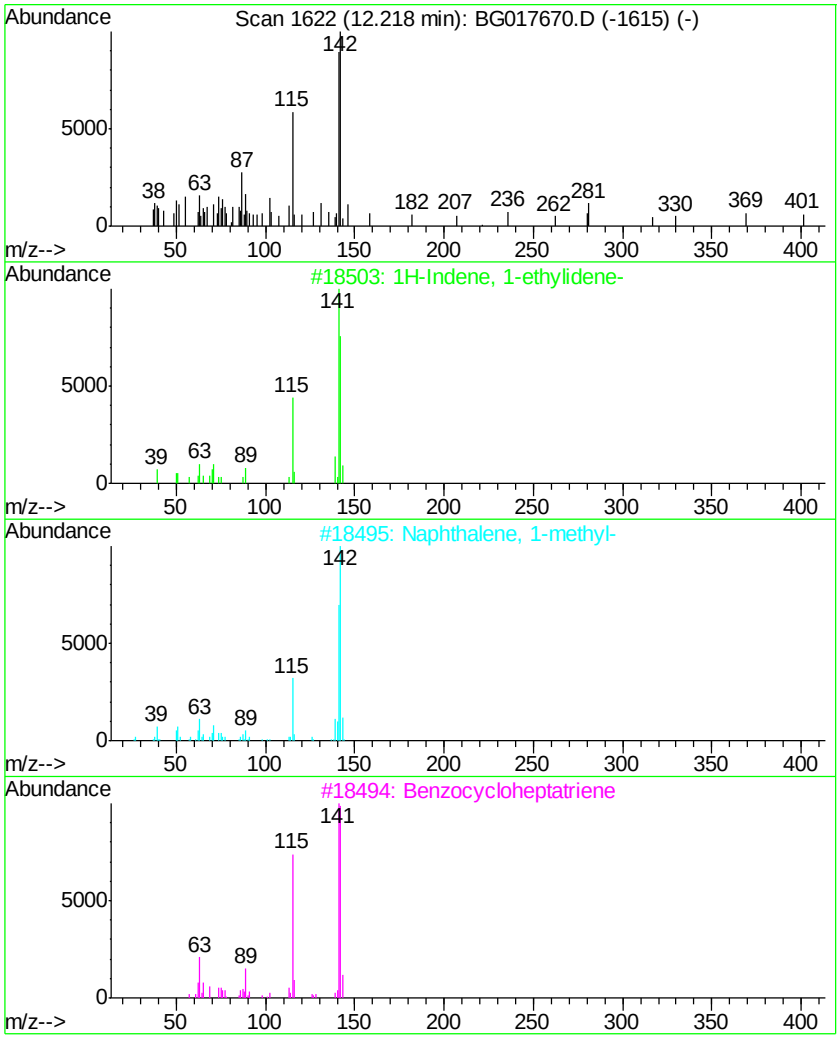
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

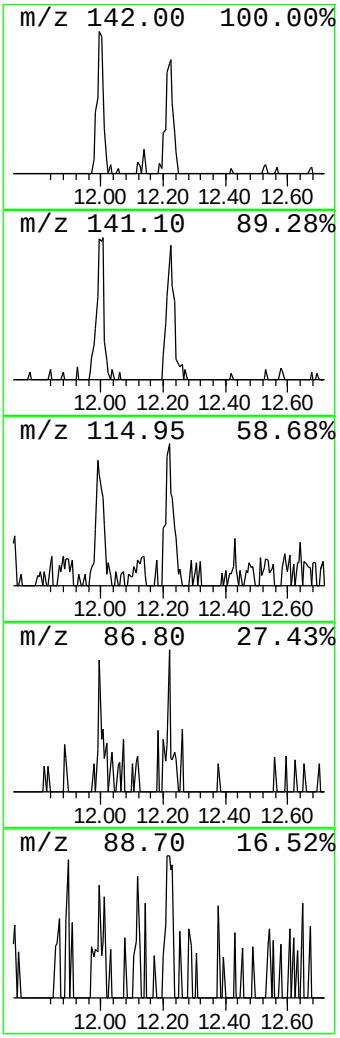
****

Peak Number 9 1H-Indene, 1-ethylidene- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.22	4.40 ng	37679	Naphthalene-d8	10.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	64
3	Benzocycloheptatriene	142	C11H10	000264-09-5	64
4	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	47
5	3,5-Difluorobenzaldehyde	142	C7H4F2O	032085-88-4	47





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

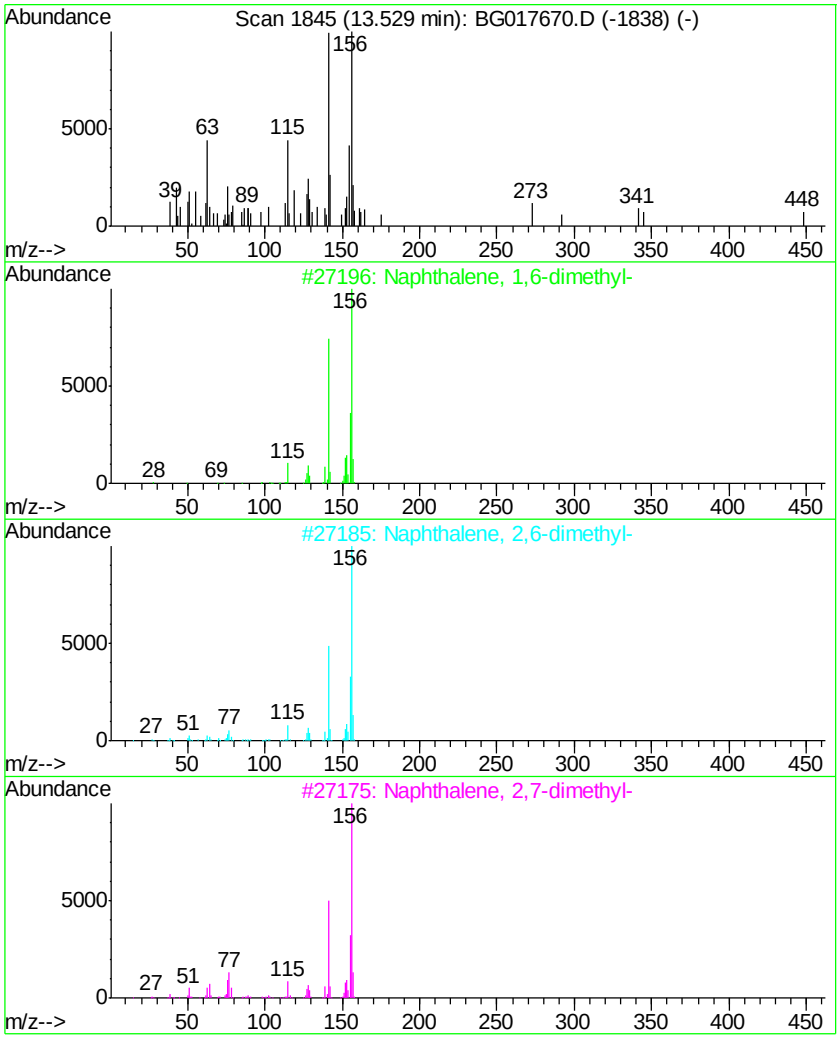
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

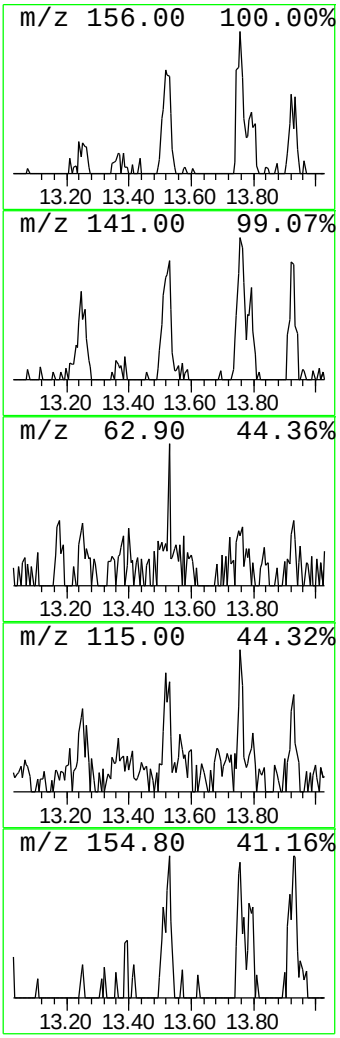
****

Peak Number 10 Naphthalene, 1,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	2.83 ng	40800	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	81
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	46
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	46
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	46
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	43





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

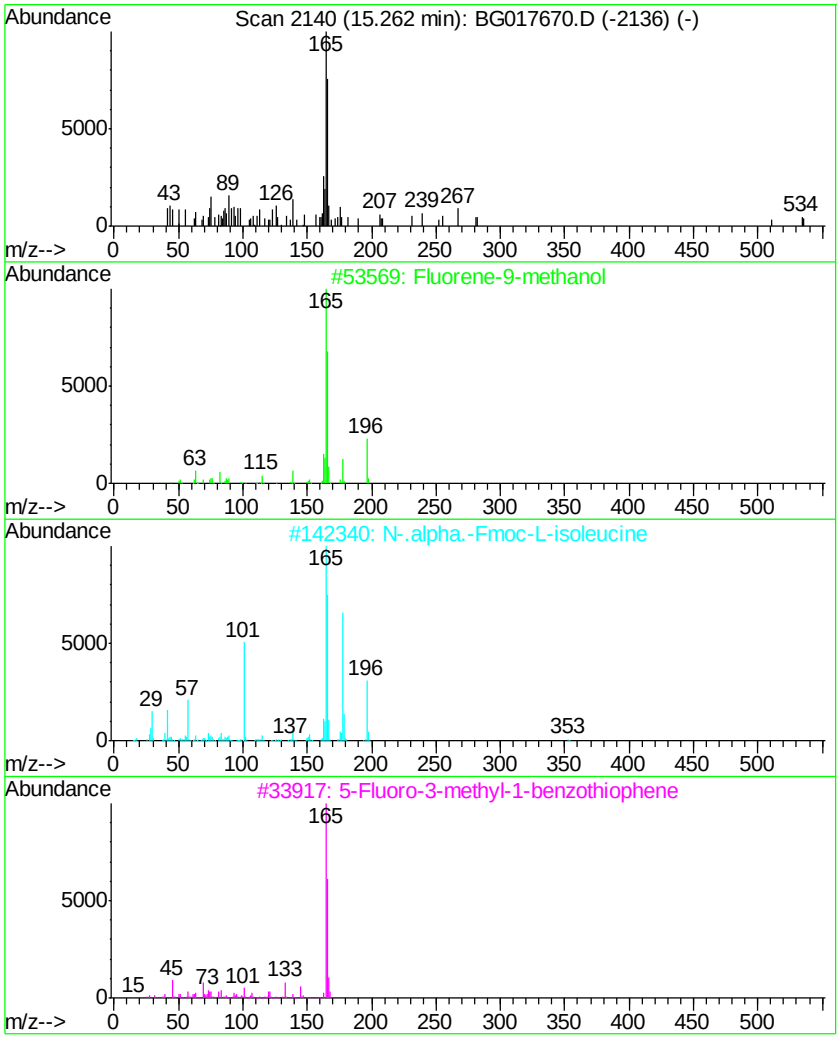
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

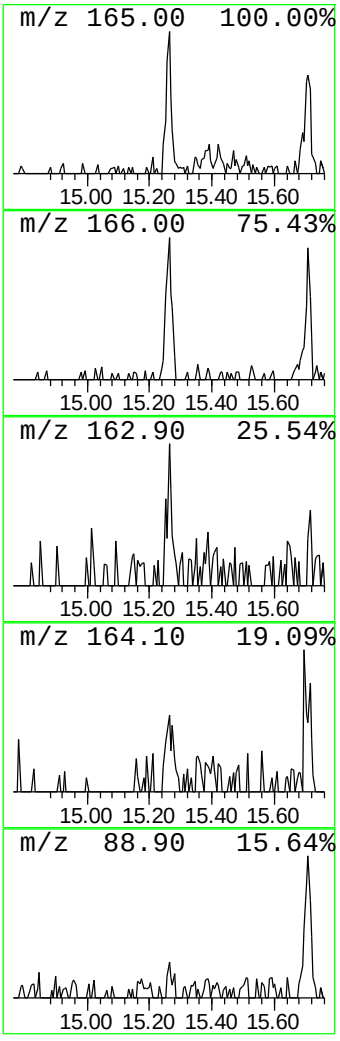
****

Peak Number 11 Fluorene-9-methanol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	2.23 ng	32190	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluorene-9-methanol	196	C14H12O	024324-17-2	64
2	N-.alpha.-Fmoc-L-isoleucine	353	C21H23NO4	071989-23-6	59
3	5-Fluoro-3-methyl-1-benzothiophene	166	C9H7FS	1000298-17-0	58
4	Benzaldehyde, 2,4-dihydroxy-3,6-...	166	C9H10O3	034883-14-2	53
5	Benzene, 1,1'-(diazomethylene)bis-	194	C13H10N2	000883-40-9	53





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

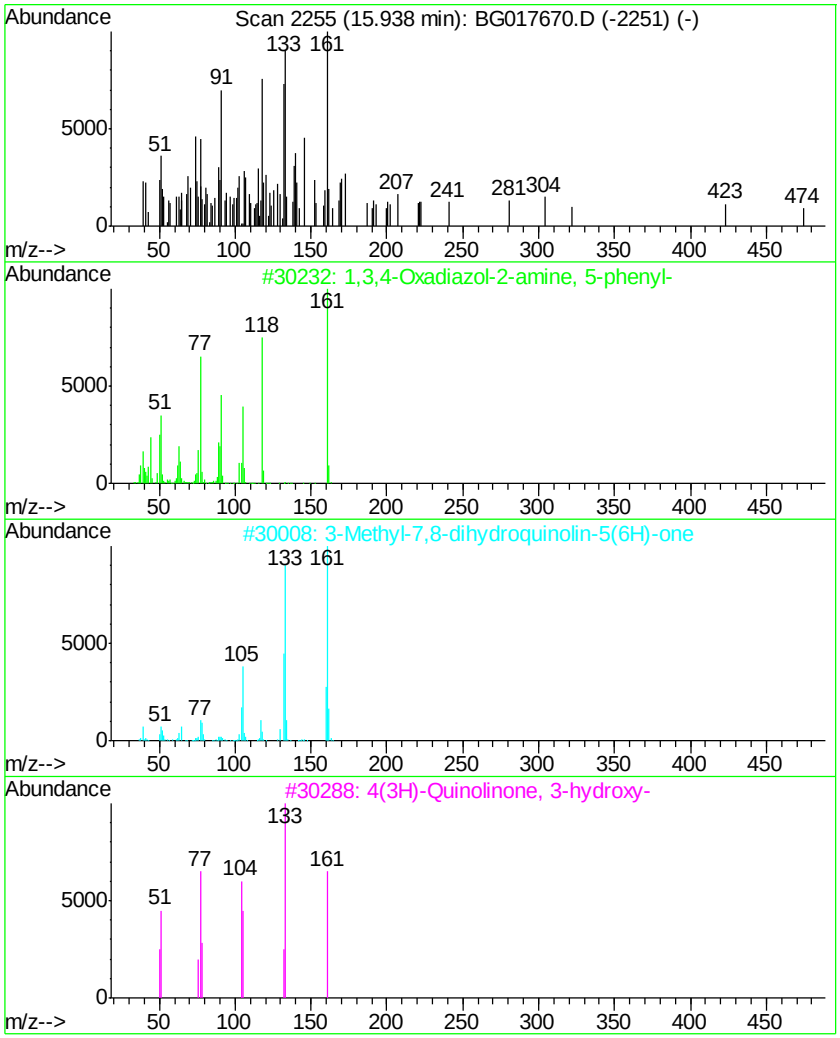
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

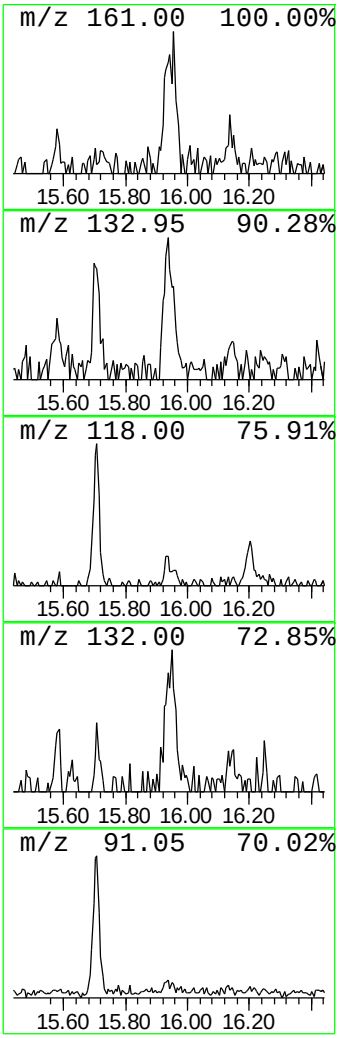
****

Peak Number 12 unknown15.94 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	2.58 ng	55282	Phenanthrene-d10	16.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,4-Oxadiazol-2-amine, 5-phenyl-	161	C8H7N3O	001612-76-6	42
2	3-Methyl-7,8-dihydroquinolin-5(6...	161	C10H11NO	060247-70-3	38
3	4(3H)-Quinolinone, 3-hydroxy-	161	C9H7NO2	055759-82-5	35
4	3-Phenyl-.delta.-1,2,4-triazolin...	161	C8H7N3O	000939-07-1	30
5	5-Chloro-2-thiophenecarbaldehyde...	161	C5H4ClNOS	014345-95-0	27





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

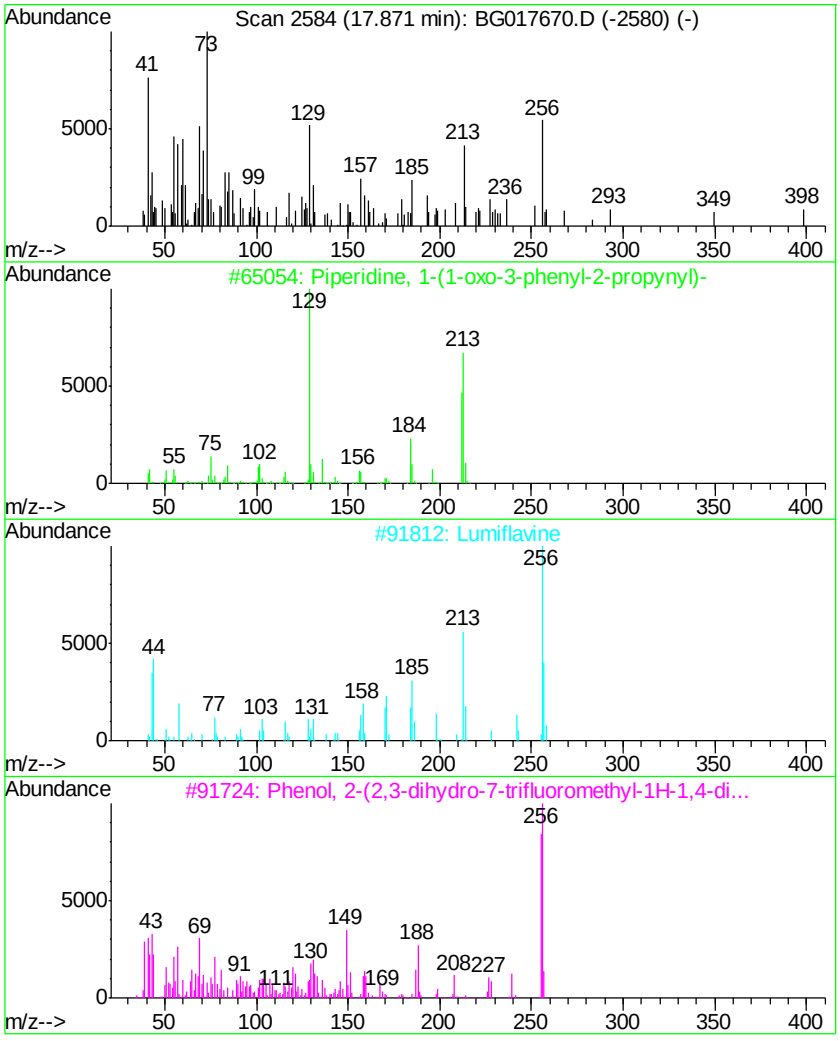
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

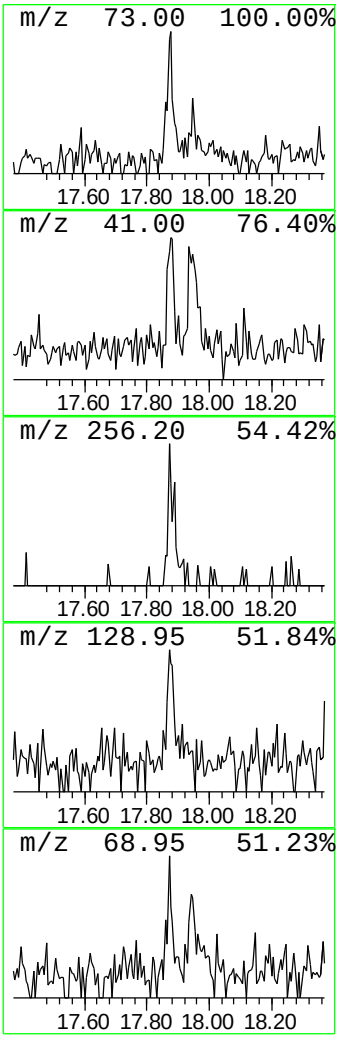
****

Peak Number 13 unknown17.87 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	2.28 ng	48974	Phenanthrene-d10	16.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Piperidine, 1-(1-oxo-3-phenyl-2-...	213	C14H15NO	014143-92-1	30
2	Lumiflavine	256	C13H12N4O2	001088-56-8	27
3	Phenol, 2-(2,3-dihydro-7-trifluo...	256	C12H11F3N2O	240417-93-0	11
4	p-(4-Ethylcyclohexyl)benzonitrile	213	C15H19N	073592-81-1	10
5	2-Oxabicyclo[4.2.0]oct-4-en-3-on...	512	C32H32O6	1000159-98-4	10





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

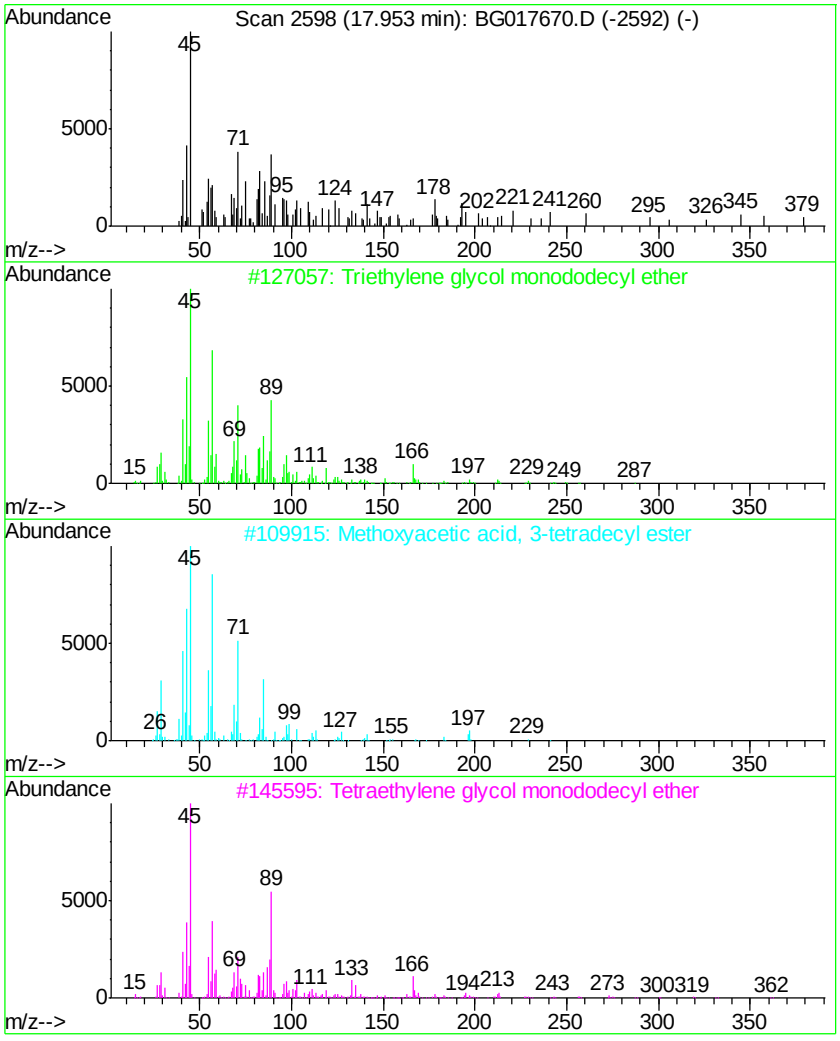
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

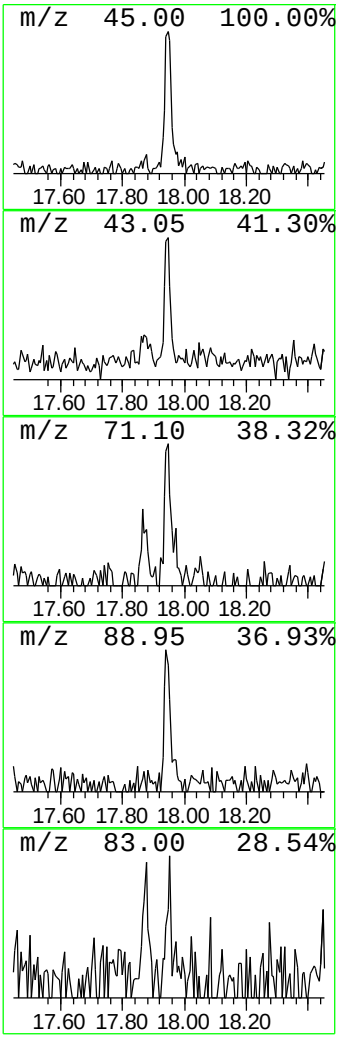
****

Peak Number 14 unknown17.95 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	3.38 ng	72537	Phenanthrene-d10	16.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triethylene glycol monododecyl e...	318	C18H38O4	003055-94-5	43
2	Methoxyacetic acid, 3-tetradecyl...	286	C17H34O3	1000282-04-9	43
3	Tetraethylene glycol monododecyl...	362	C20H42O5	005274-68-0	32
4	Methoxyacetic acid, 2-butyl ester	146	C7H14O3	070159-90-9	27
5	2-Tetradecanol	214	C14H30O	004706-81-4	22





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061215\
Data File : BG017670.D
Acq On : 13 Jun 2015 11:52
Operator : TP/IZ
Sample : G2627-12
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CMW-37

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061215.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
unknown7.07	7.07	88.5	ng	635486	1	7.52	143635	20.0
Indane	7.89	10.7	ng	77068	1	7.52	143635	20.0
Phenol, 2,6-dimet...	8.96	4.3	ng	37125	2	10.31	171272	20.0
Benzene, 1-etheny...	9.72	9.4	ng	80525	2	10.31	171272	20.0
Phenol, 3,5-dimet...	9.84	9.6	ng	82441	2	10.31	171272	20.0
unknown11.43	11.43	3.8	ng	32853	2	10.31	171272	20.0
unknown11.89	11.89	3.7	ng	31604	2	10.31	171272	20.0
1H-Indene, 1-ethy...	12.22	4.4	ng	37679	2	10.31	171272	20.0
Naphthalene, 1,6-...	13.53	2.8	ng	40800	3	14.20	288393	20.0
Fluorene-9-methanol	15.26	2.2	ng	32190	3	14.20	288393	20.0
unknown15.94	15.94	2.6	ng	55282	4	16.96	429044	20.0
unknown17.87	17.87	2.3	ng	48974	4	16.96	429044	20.0
unknown17.95	17.95	3.4	ng	72537	4	16.96	429044	20.0