LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

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-BG122116.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	5.259	407	412	420	rBV	206902	342624	4.05%	0.953%
2	5.741	487	494	503	rBV	646289	1176552	13.91%	3.272%
3	5.817	503	507	516	rVB3	68750	127751	1.51%	0.355%
4	7.363	764	770	784	rVB	416283	846755	10.01%	2.355%
5	7.568	798	805	812	rBV3	52040	99845	1.18%	0.278%
6	7.733	826	833	845	rVB	1196819	2187236	25.86%	6.083%
7	8.197	905	912	921	rBV	234476	445750	5.27%	1.240%
8	8.303	921	930	939	rVB	320151	583960	6.90%	1.624%
9	8.526	961	968	981	rVB	1083661	2051546	24.26%	5.705%
10	8.673	987	993	1000	rVB2	114788	200621	2.37%	0.558%
11	8.826	1012	1019	1024	rBV2	58896	111891	1.32%	0.311%
12	9.296	1091	1099	1106	rBV3	70062	155449	1.84%	0.432%
13	9.384	1106	1114	1126	rVV	1207508	2257381	26.69%	6.278%
14	9.825	1182	1189	1194	rBV2	109934	193029	2.28%	0.537%
15	10.253	1256	1262	1271	rVB	156038	307955	3.64%	0.856%
16	10.406	1279	1288	1302	rBV	307924	626337	7.41%	1.742%
17	11.029	1389	1394	1399	rBV2	284274	528078	6.24%	1.469%
18	11.076	1399	1402	1408	rVB	216339	355035	4.20%	0.987%
19	12.045	1559	1567	1571	rBV6	43178	86389	1.02%	0.240%
20	12.239	1592	1600	1609	rBV5	72485	142539	1.69%	0.396%
21	12.886	1703	1710	1718	rVB4	87793	173642	2.05%	0.483%
22	13.450	1798	1806	1815	rVB	2643620	4161786	49.21%	11.574%
23	13.808	1860	1867	1873	rBV3	80699	139897	1.65%	0.389%
24	13.973	1890	1895	1900	rBV3	45347	90679	1.07%	0.252%
25	14.266	1940	1945	1950	rBV2	77305	120229	1.42%	0.334%
26	14.825	2031	2040	2052	rBV	568307	955967	11.30%	2.658%
27	16.317	2287	2294	2305	rBV	2672212	4295974	50.79%	11.947%
28	17.574	2502	2508	2516	rBV	735724	1176696	13.91%	3.272%
29	18.409	2645	2650	2658	rBV5	58777	88948	1.05%	0.247%
30	20.154	2940	2947	2953	rBV	6408869	8457768	100.00%	23.520%
31	21.863	3232	3238	3245	rVB	982231	1705373	20.16%	4.743%
32	25.259	3806	3816	3830	rVB4	561816	1765561	20.88%	4.910%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

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-BG122116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

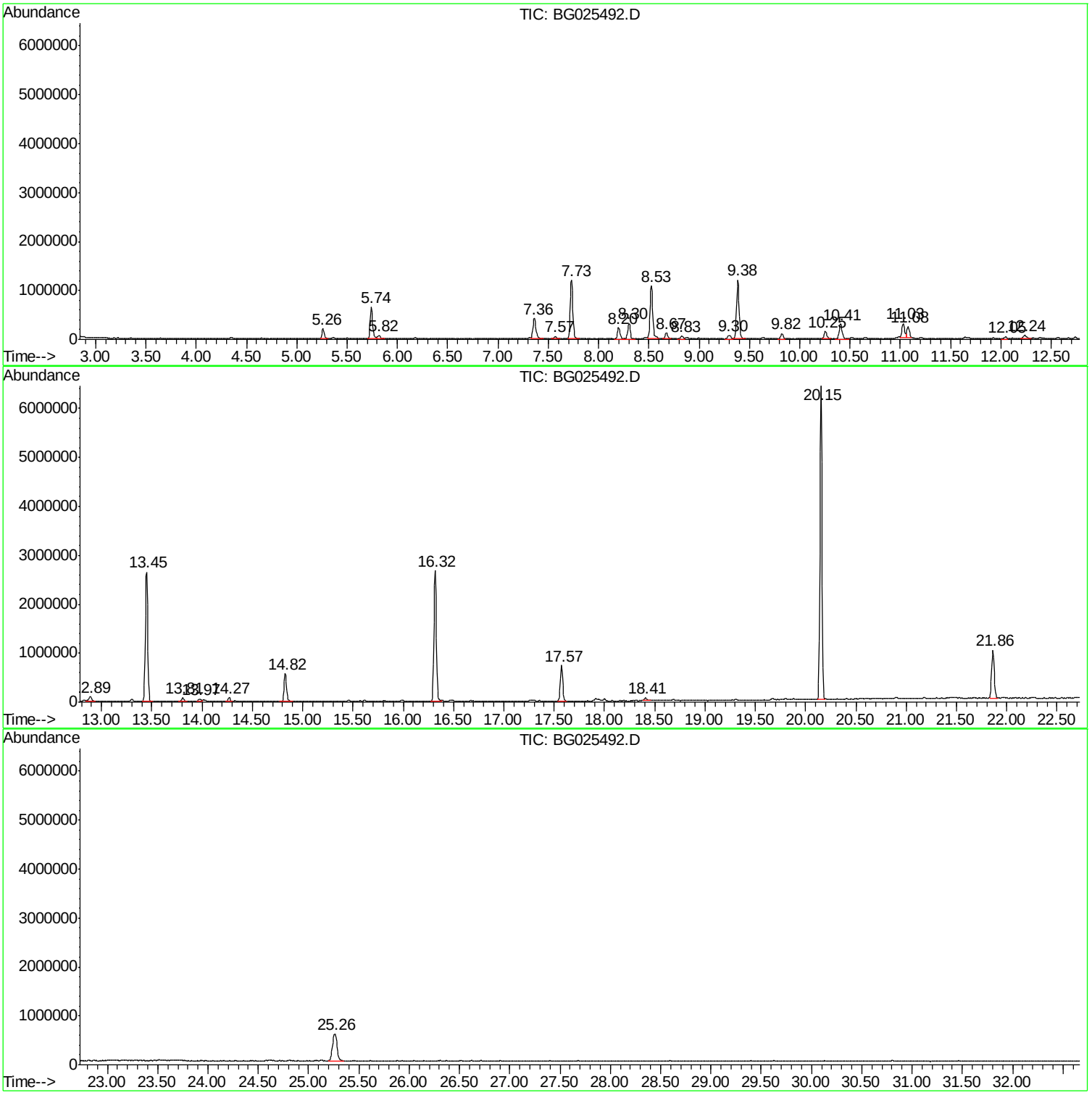
Sum of corrected areas: 35959243

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

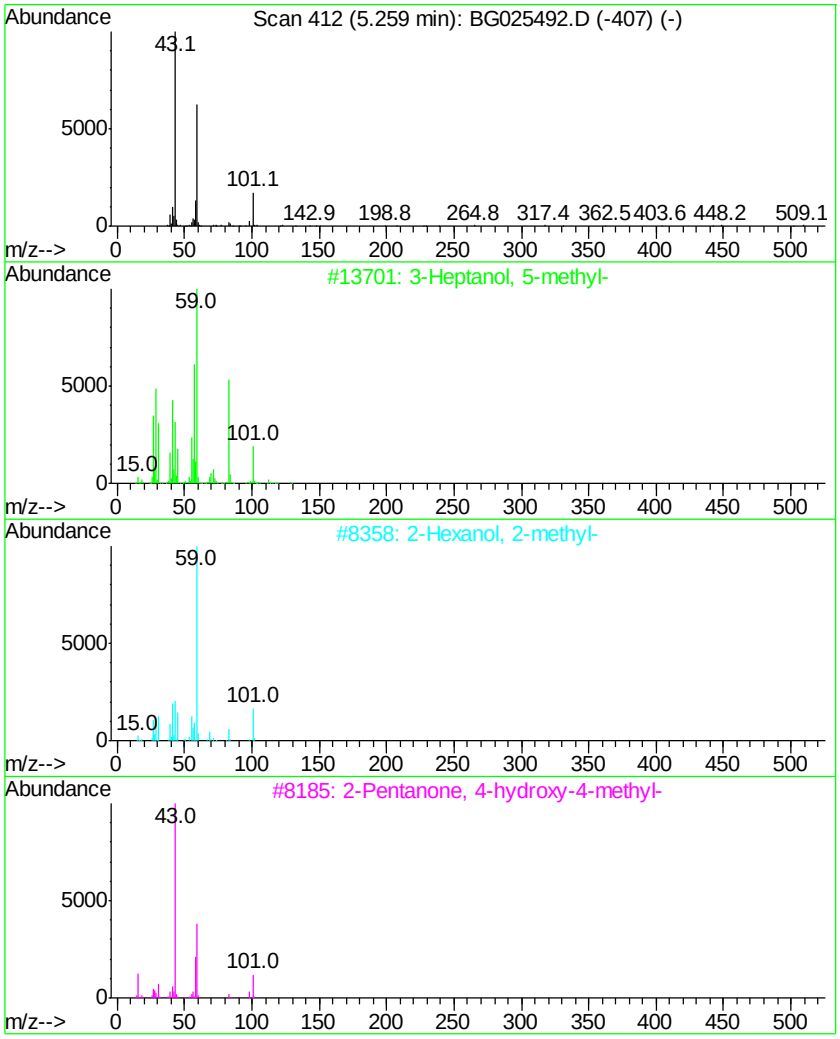
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

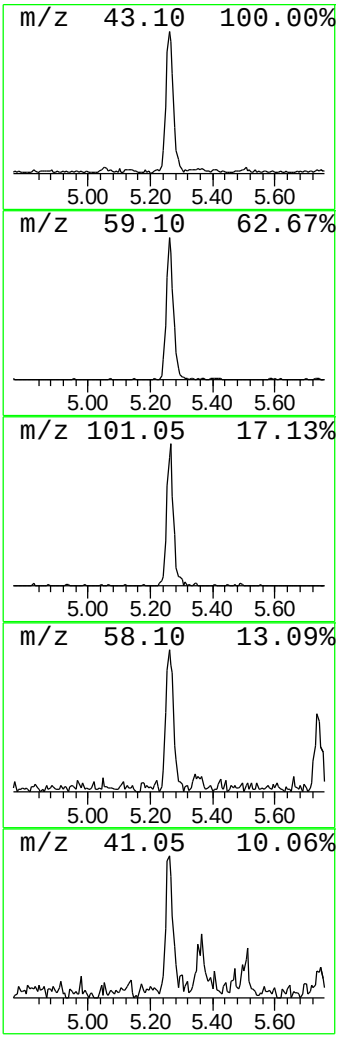
****

Peak Number 1 unknown5.26 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.26	15.37 ng	342624	1,4-Dichlorobenzene-d4	8.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptanol, 5-methyl-	130	C8H18O	018720-65-5	28
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	23
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

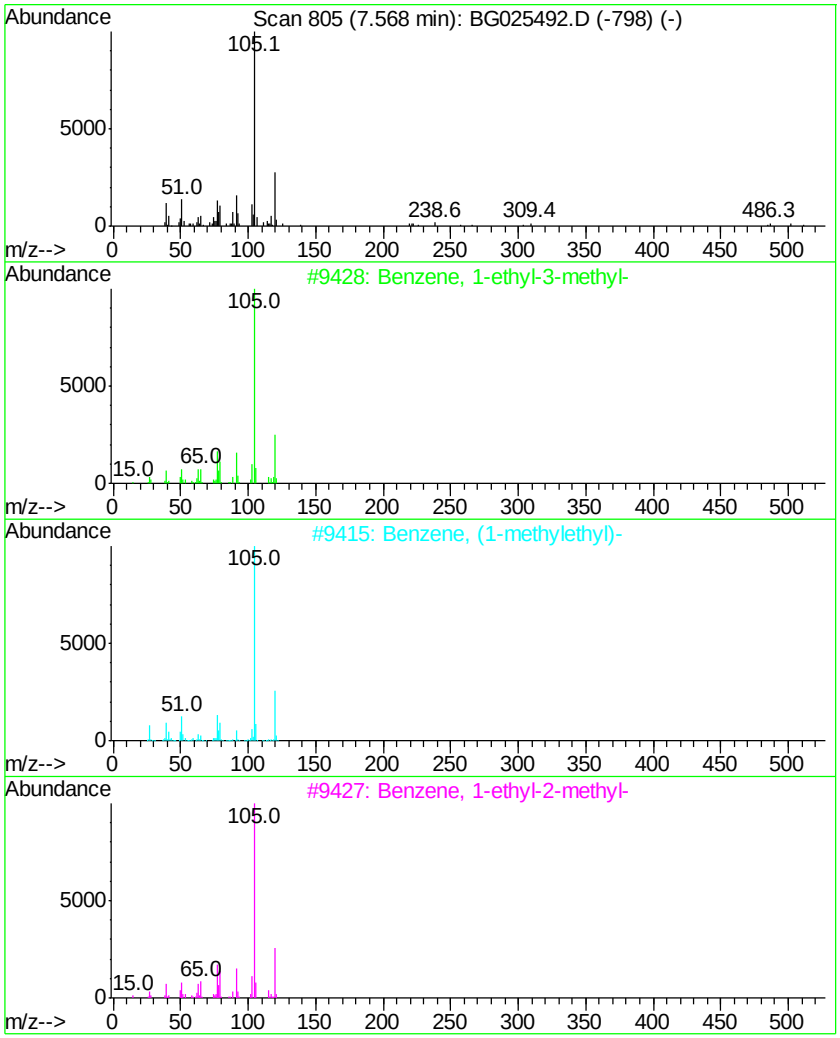
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

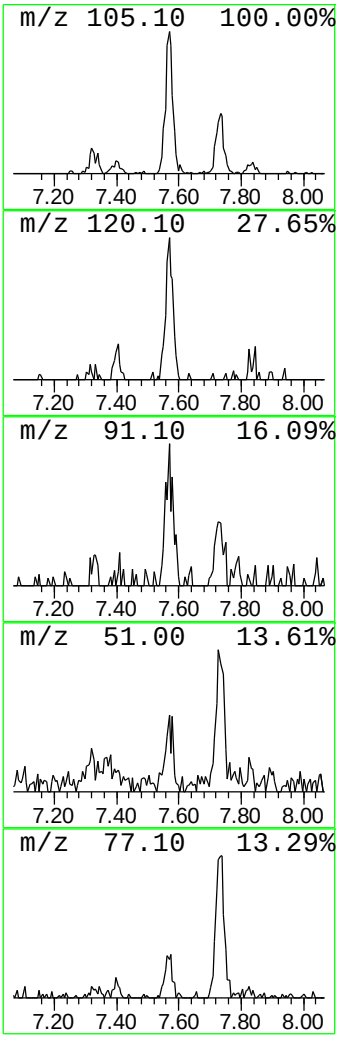
****

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	4.48 ng	99845	1,4-Dichlorobenzene-d4	8.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	68
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	64
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	64





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

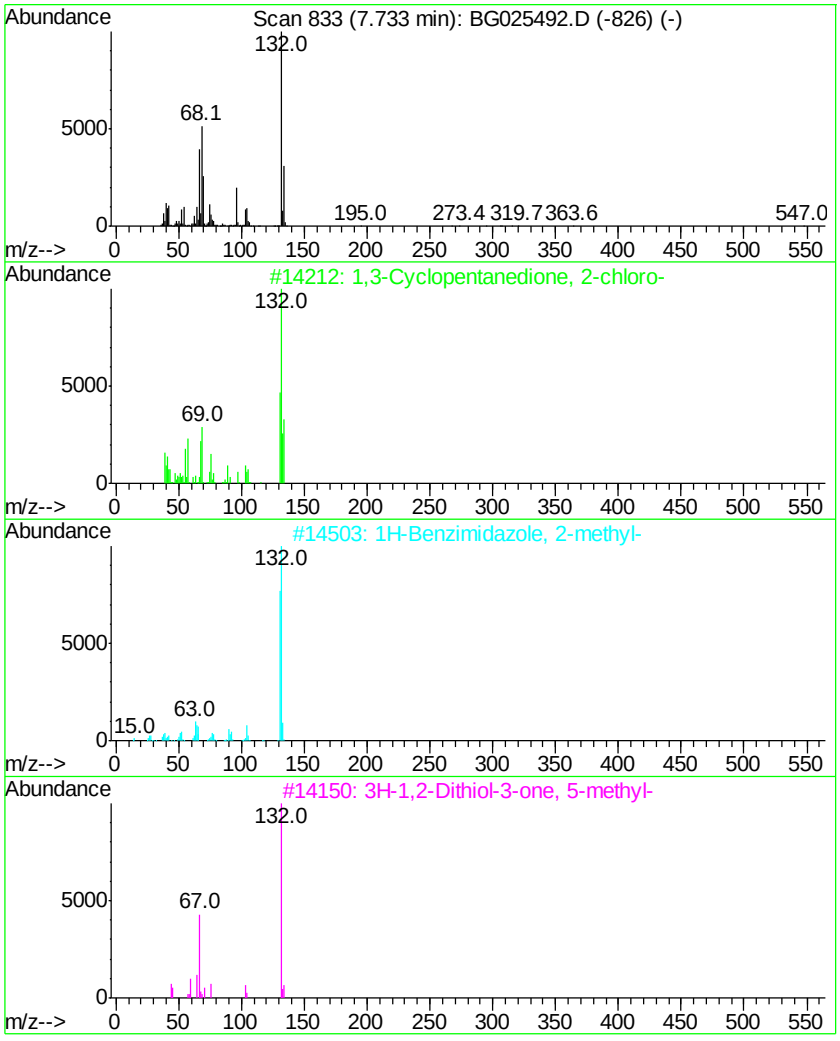
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

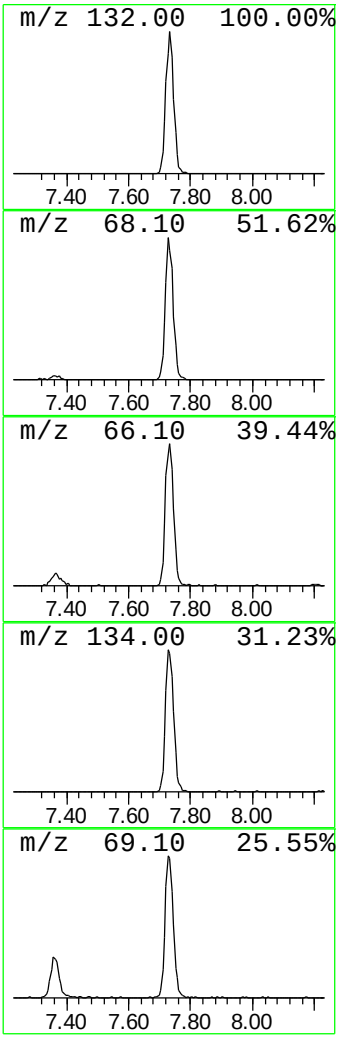
****

Peak Number 4 unknown7.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	98.14 ng	2187240	1,4-Dichlorobenzene-d4	8.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	27
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
3	3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
4	1H-Indenol	132	C9H8O	056631-57-3	10
5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	10





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

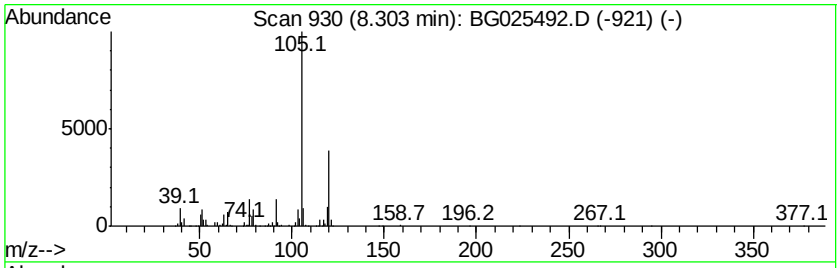
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

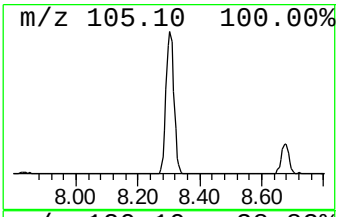
****

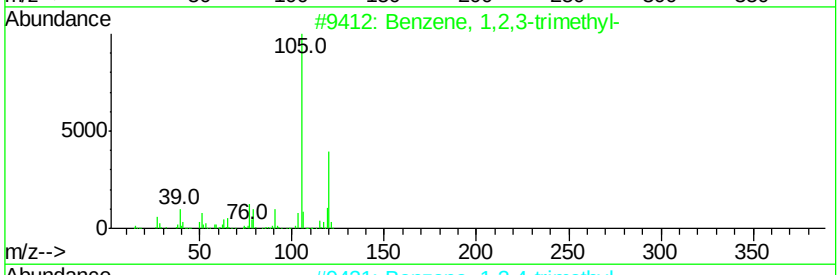
Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 3

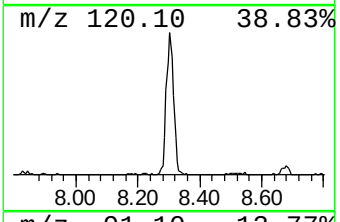
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.30	26.20 ng	583960	1,4-Dichlorobenzene-d4	8.20

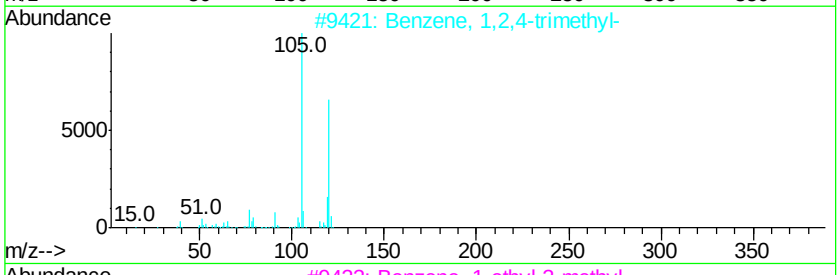
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90

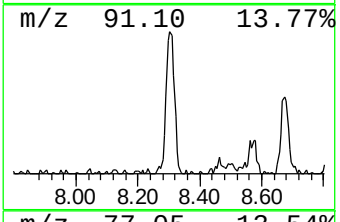


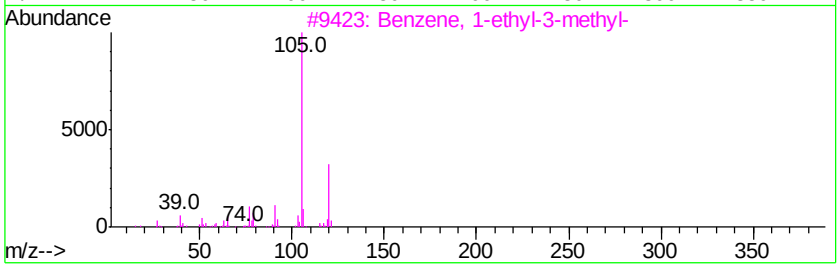


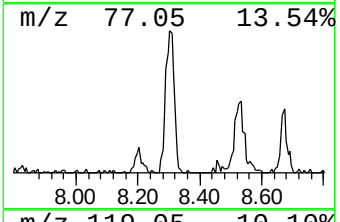


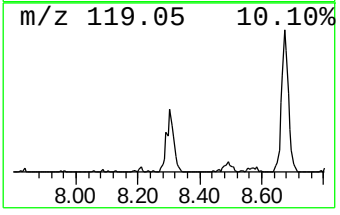












Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

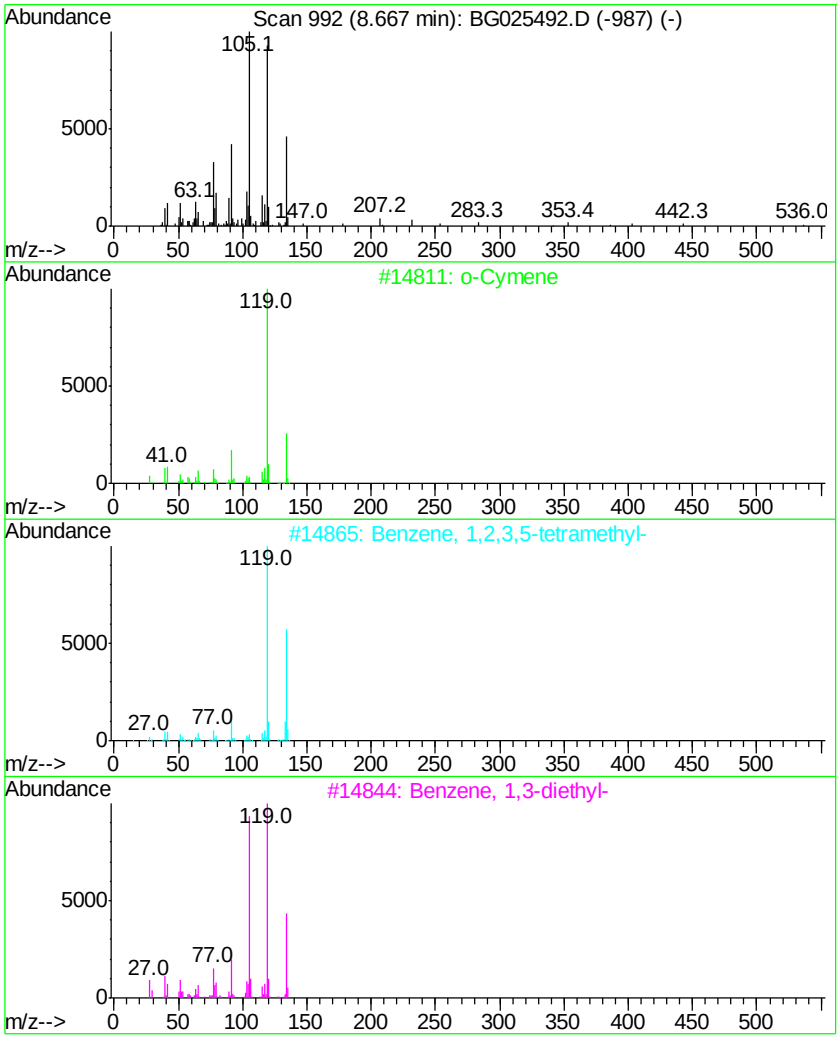
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

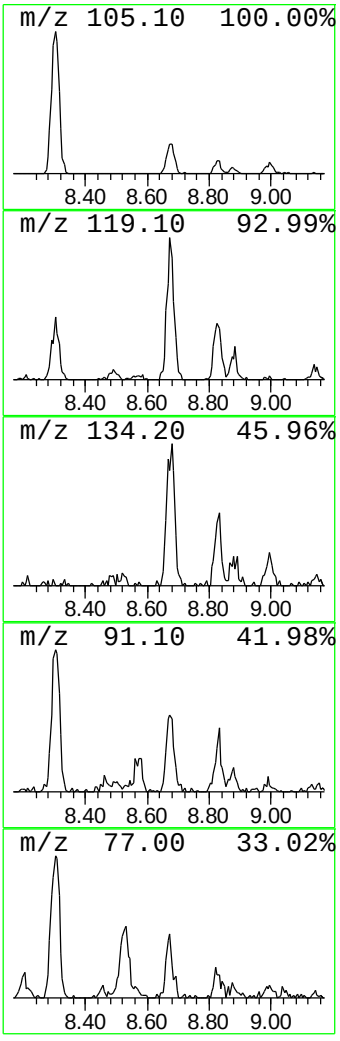
****

Peak Number 7 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.67	9.00 ng	200621	1,4-Dichlorobenzene-d4	8.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	76
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	64
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	64
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	64
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

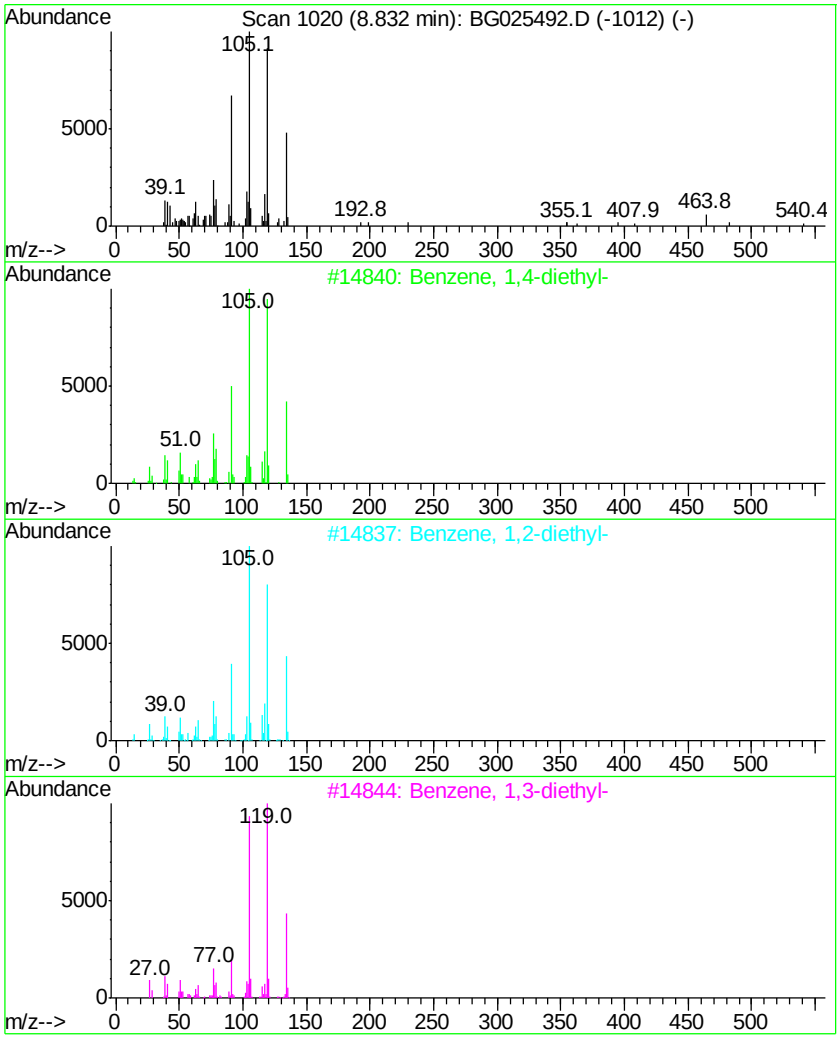
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

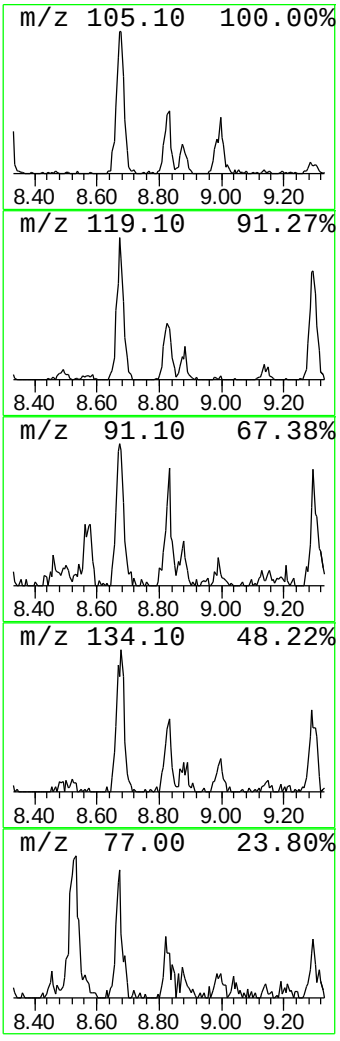
****

Peak Number 8 Benzene, 1,4-diethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.83	5.02 ng	111891	1,4-Dichlorobenzene-d4	8.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	74
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	72
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	64
4	5H-5-Methyl-6,7-dihydrocyclopent...	134	C8H10N2	023747-48-0	47
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	35





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

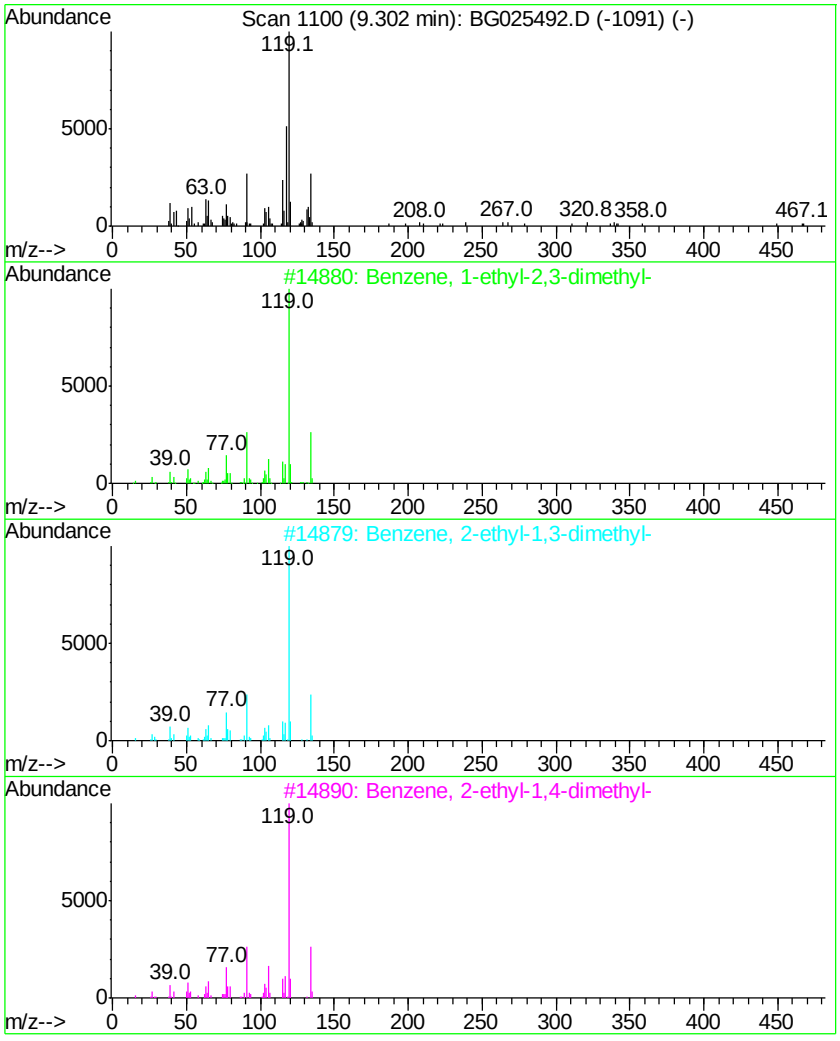
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

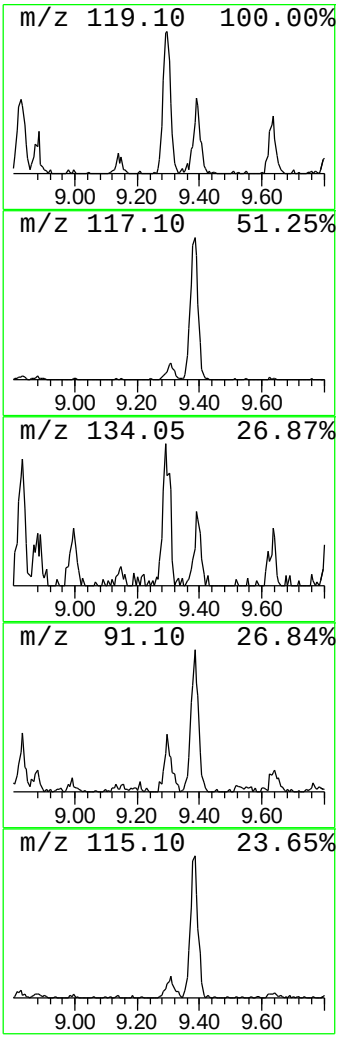
****

Peak Number 9 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.30	6.97 ng	155449	1,4-Dichlorobenzene-d4	8.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	76
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	68
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	68
5	o-Cymene	134	C10H14	000527-84-4	64





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

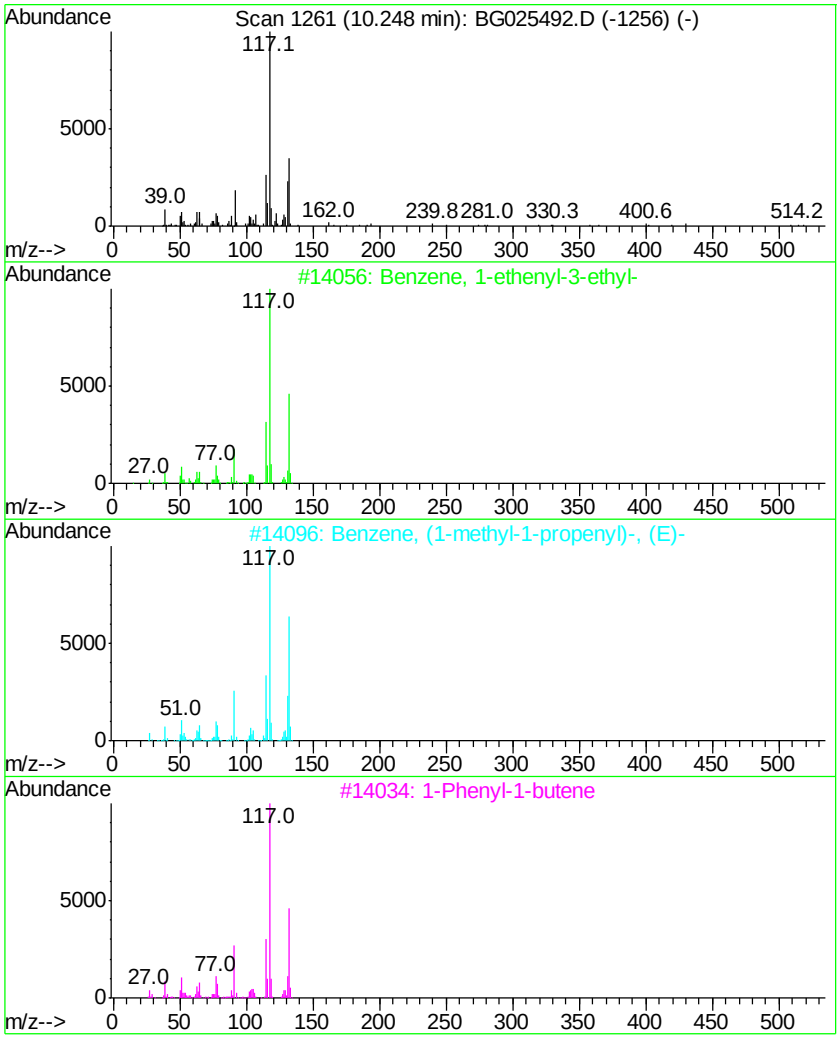
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

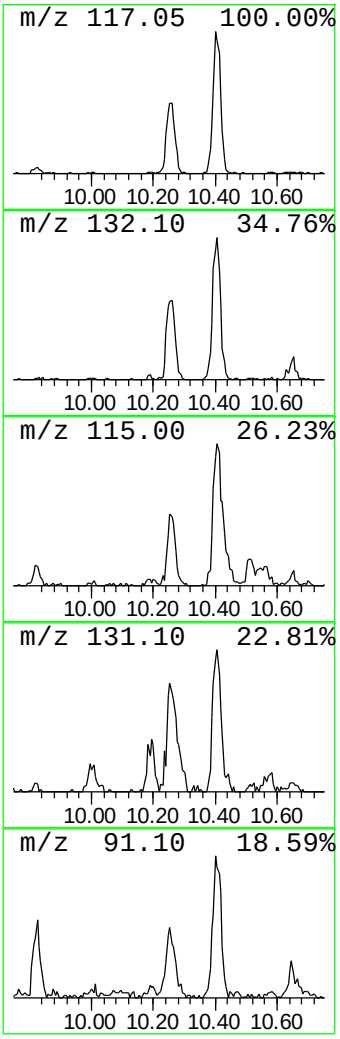
****

Peak Number 11 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	11.66 ng	307955	Naphthalene-d8	11.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	93
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.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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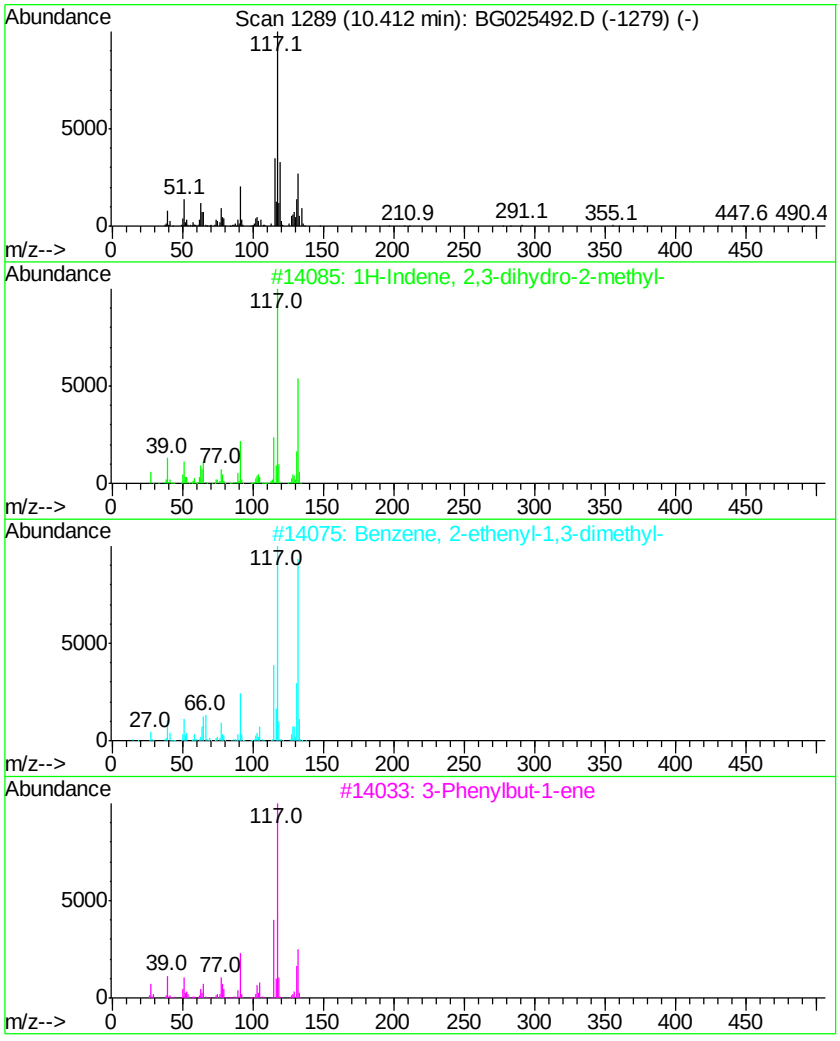
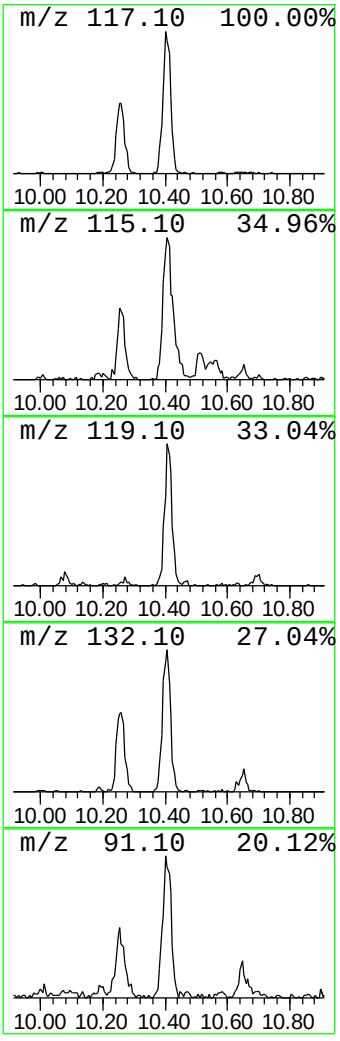
Peak Number 12 1H-Indene, 2,3-dihydro-2-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.41	23.72 ng	626337	Naphthalene-d8	11.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	76
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	76
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	76
5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	70





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.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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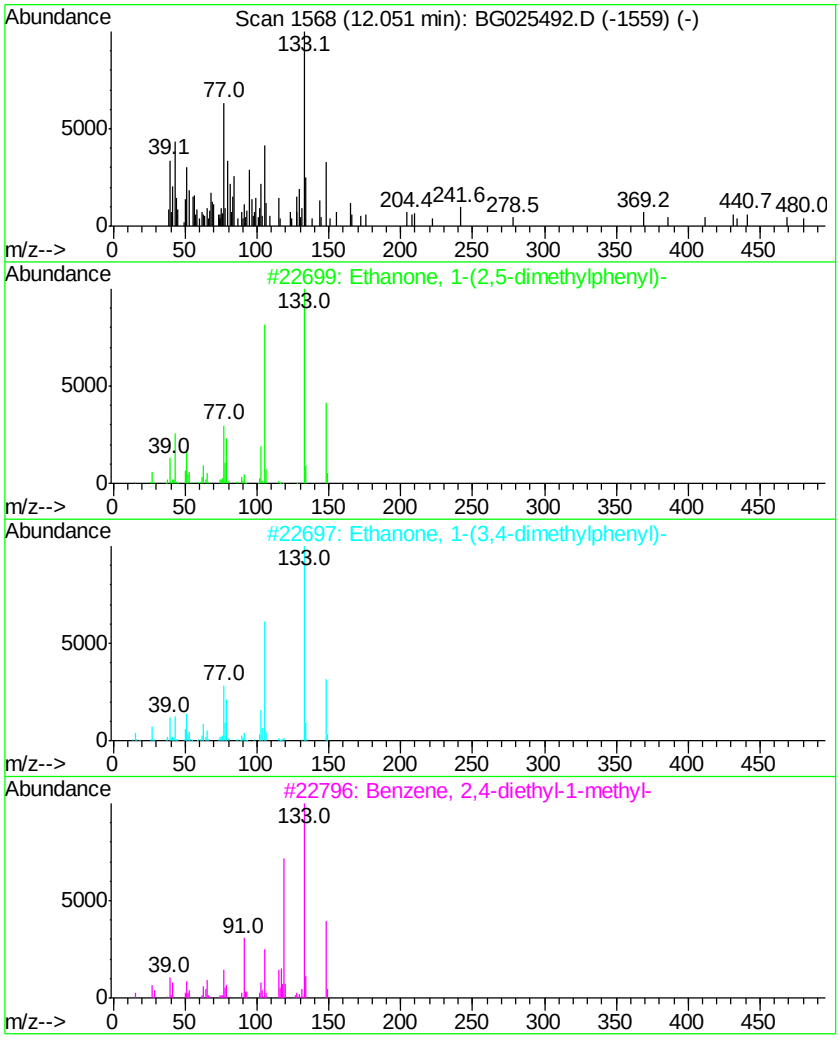
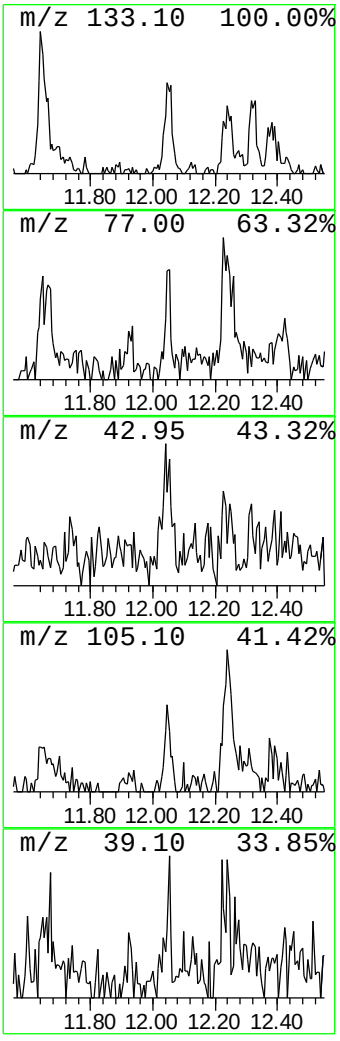
Peak Number 13 unknown12.05 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	3.27 ng	86389	Naphthalene-d8	11.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	49
2	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	49
3	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	47
4	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	47
5	Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	47





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.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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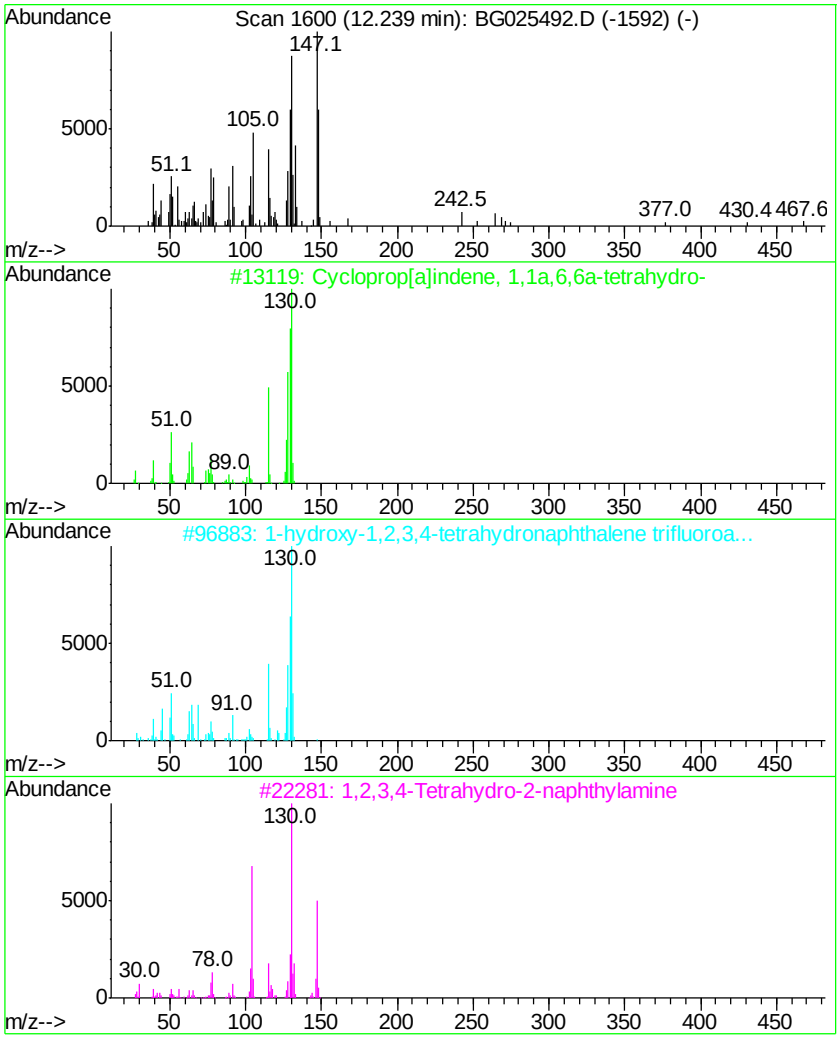
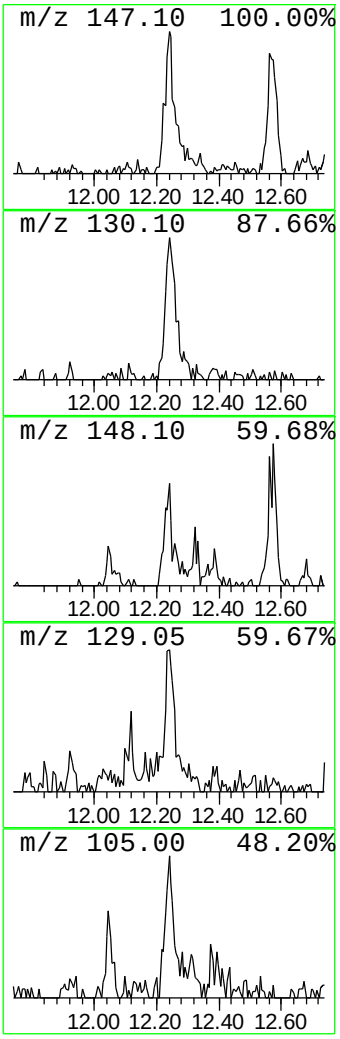
Peak Number 14 unknown12.24 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.24	5.40 ng	142539	Naphthalene-d8	11.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	38
2	1-hydroxy-1,2,3,4-tetrahydronaph...	244	C12H11F3O2	1000376-44-1	38
3	1,2,3,4-Tetrahydro-2-naphthylamine	147	C10H13N	002954-50-9	38
4	Benzene, 1,3-butadienyl-	130	C10H10	001515-78-2	30
5	3-Methylbenzothiophene	148	C9H8S	001455-18-1	25





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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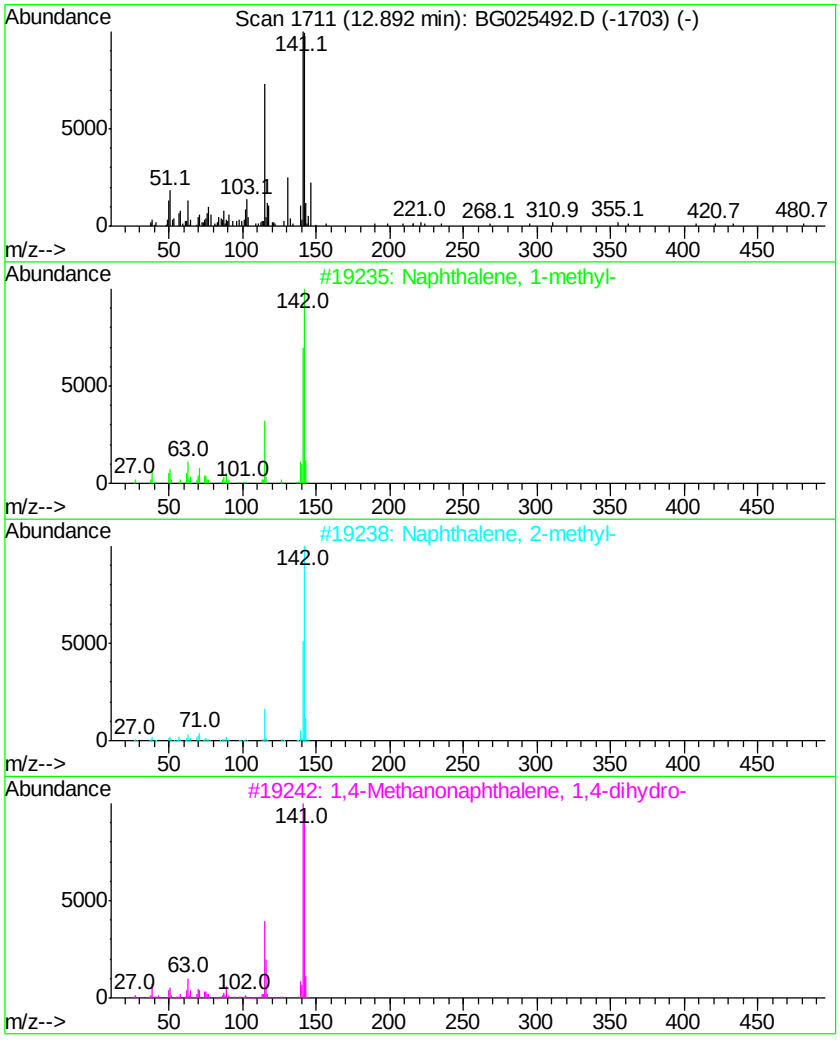
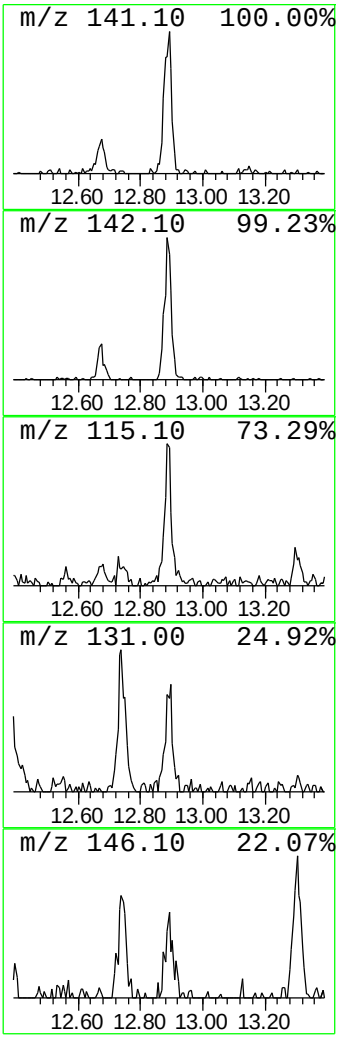
Peak Number 15 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	6.58 ng	173642	Naphthalene-d8	11.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	68
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	58
4	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	49
5	5-Methylisophthalonitrile	142	C9H6N2	039718-07-5	49





Data Path : Z:\HPCHEM1\BNA G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.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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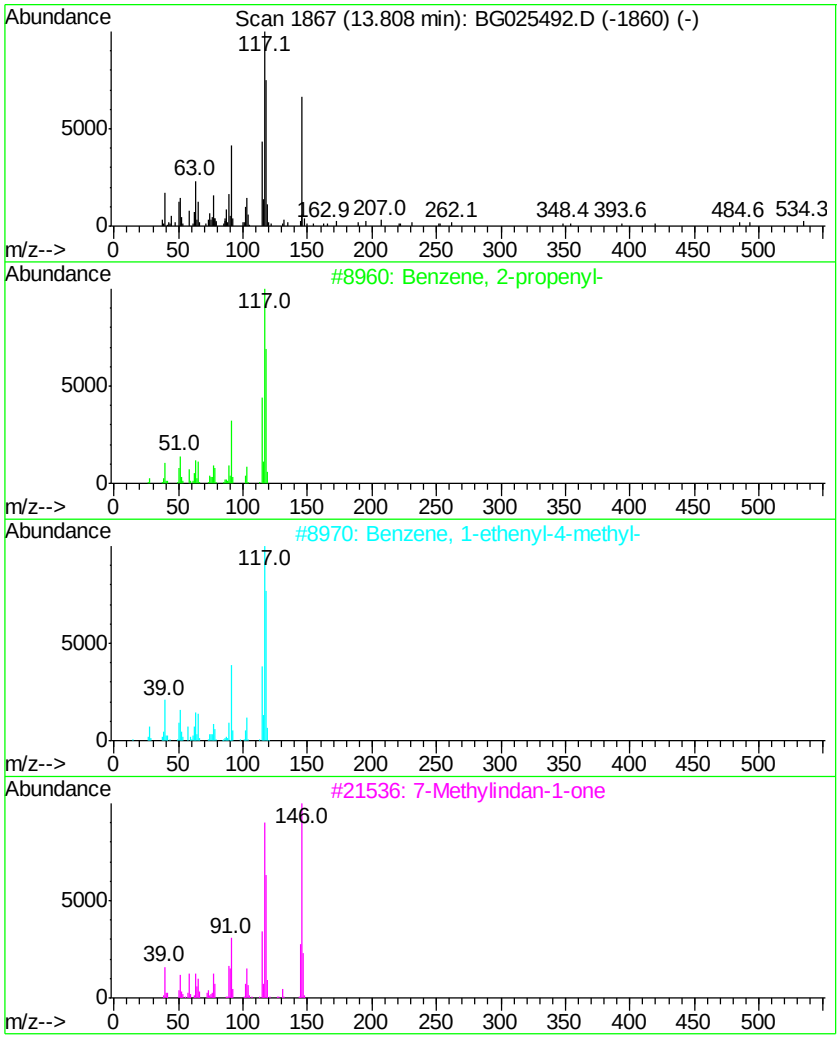
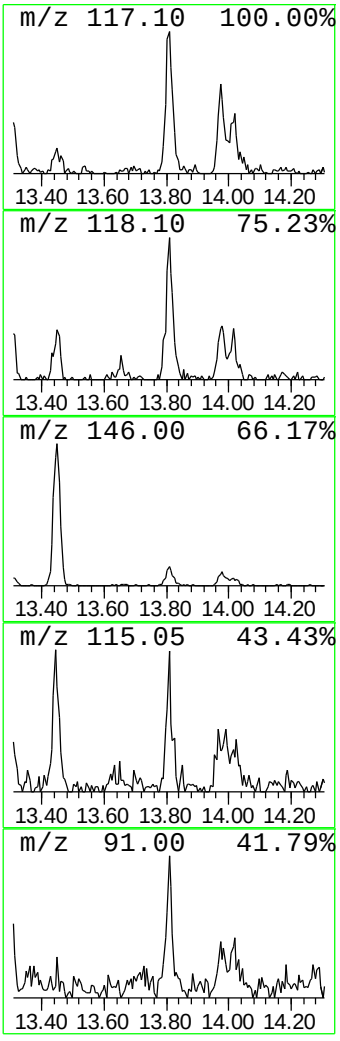
Peak Number 16 Benzene, 2-propenyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	2.93 ng	139897	Acenaphthene-d10	14.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	93
2	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	92
3	7-Methylindan-1-one	146	C10H10O	039627-61-7	91
4	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	91
5	Indane	118	C9H10	000496-11-7	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG011217\
Data File : BG025492.D
Acq On : 12 Jan 2017 21:34
Operator : UM/SJ
Sample : I1125-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : W-33

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122116.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
unknown5.26	5.26	15.4	ng	342624	1	8.20	445750	20.0
Benzene, 1-ethyl-...	7.57	4.5	ng	99845	1	8.20	445750	20.0
unknown7.73	7.73	98.1	ng	2187240	1	8.20	445750	20.0
Benzene, 1,2,3-tr...	8.30	26.2	ng	583960	1	8.20	445750	20.0
o-Cymene	8.67	9.0	ng	200621	1	8.20	445750	20.0
Benzene, 1,4-diet...	8.83	5.0	ng	111891	1	8.20	445750	20.0
Benzene, 1-ethyl-...	9.30	7.0	ng	155449	1	8.20	445750	20.0
Benzene, 1-etheny...	10.25	11.7	ng	307955	2	11.03	528078	20.0
1H-Indene, 2,3-di...	10.41	23.7	ng	626337	2	11.03	528078	20.0
unknown12.05	12.05	3.3	ng	86389	2	11.03	528078	20.0
unknown12.24	12.24	5.4	ng	142539	2	11.03	528078	20.0
Naphthalene, 1-me...	12.89	6.6	ng	173642	2	11.03	528078	20.0
Benzene, 2-propenyl-	13.81	2.9	ng	139897	3	14.82	955967	20.0