

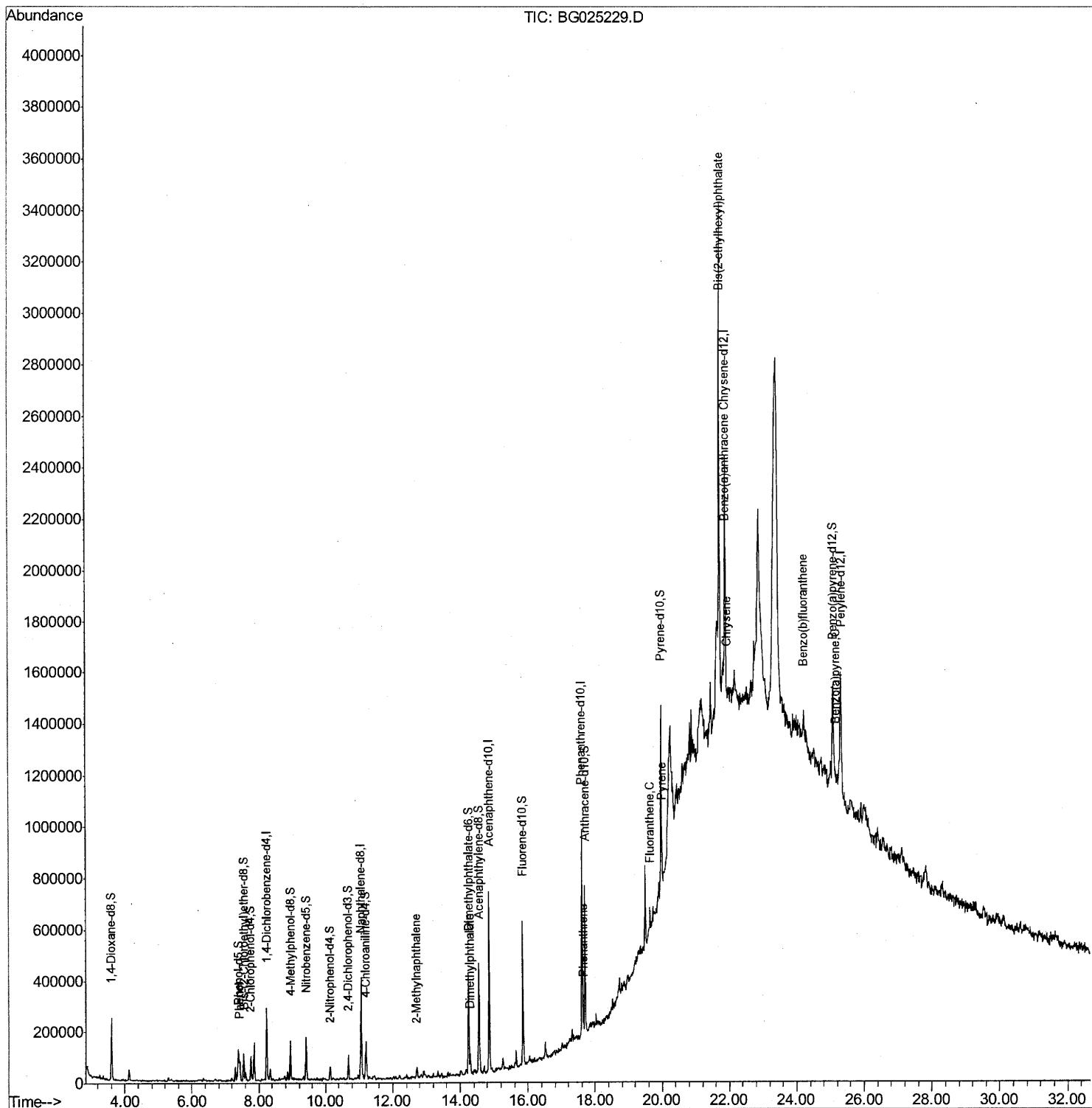
Data Path : Z:\HPCHEM1\BNA\_G\Data\BG121516\  
Data File : BG025229.D  
Acq On : 16 Dec 2016 2:29  
Operator : UM/SJ  
Sample : H5995-16  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_G  
Client Sampled :  
DA884

Manual Integrations  
APPROVED

UMANGI  
12/16/2016 6:21:22 PM

Quant Time: Dec 16 14:14:22 2016  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Dec 16 05:54:49 2016  
Response via : Initial Calibration



## Quantitation Report (Qedit)

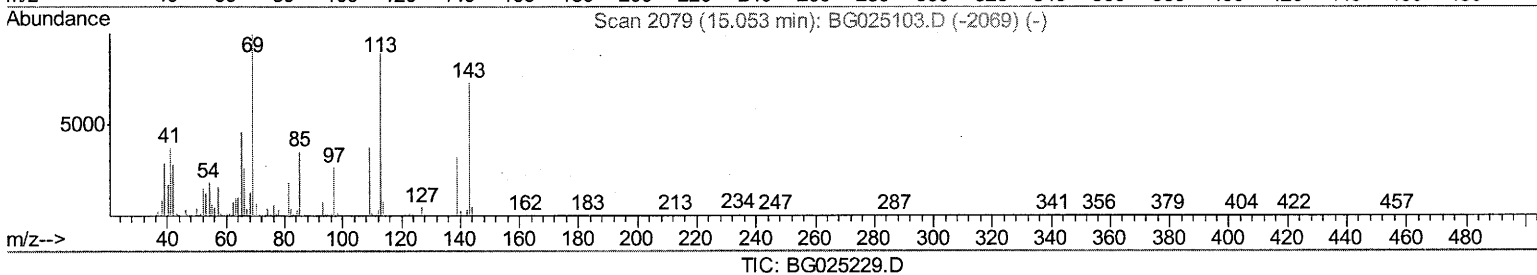
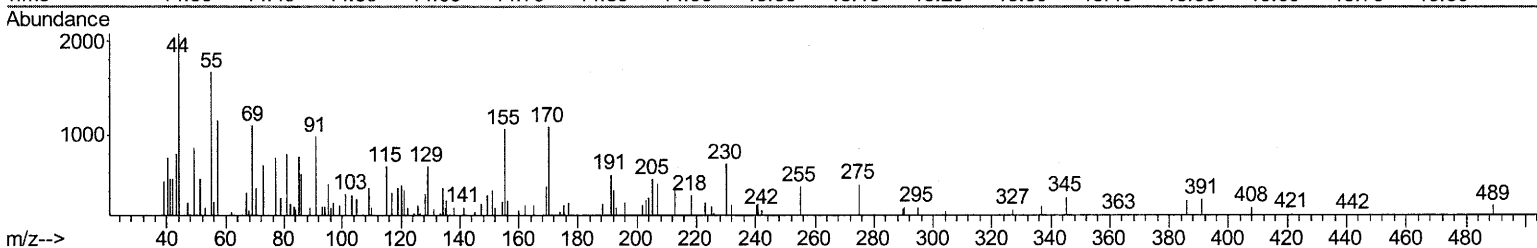
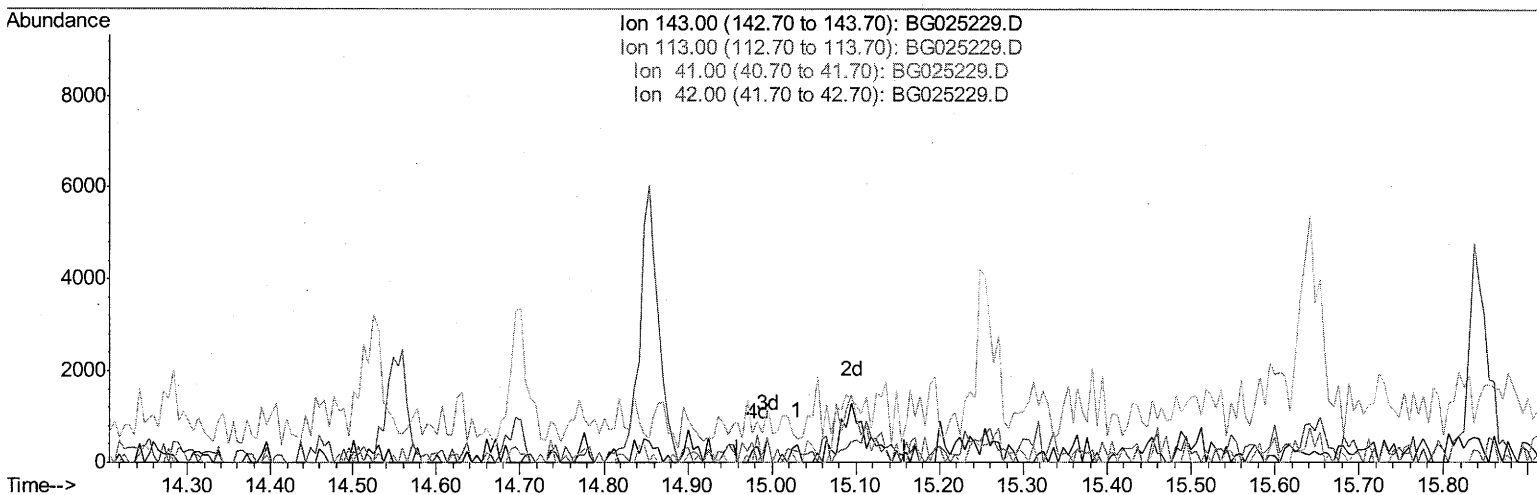
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Quant Time: Dec 16 06:20:28 2016  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Dec 16 05:54:49 2016  
Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

15.059min (-15.059) 0.00ng/ul

response 0

| Ion    | Exp%   | Act% |
|--------|--------|------|
| 143.00 | 100    | 0.00 |
| 113.00 | 117.30 | 0.00 |
| 41.00  | 58.00  | 0.00 |
| 42.00  | 42.20  | 0.00 |

# Quantitation Report (Qedit)

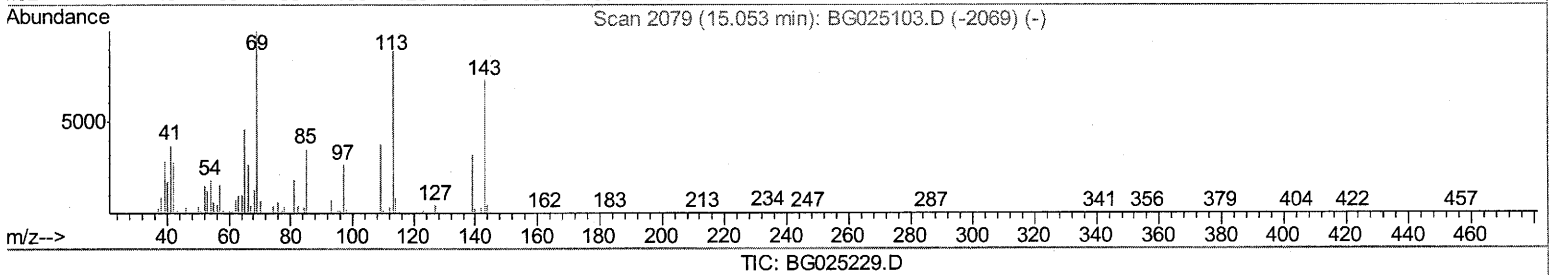
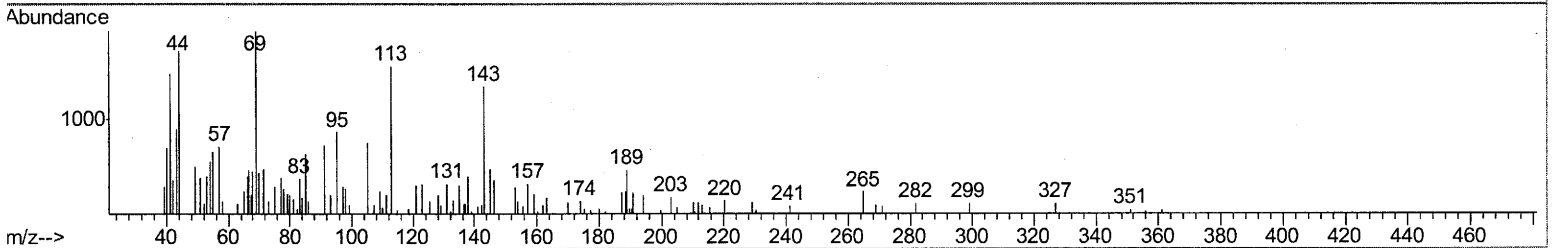
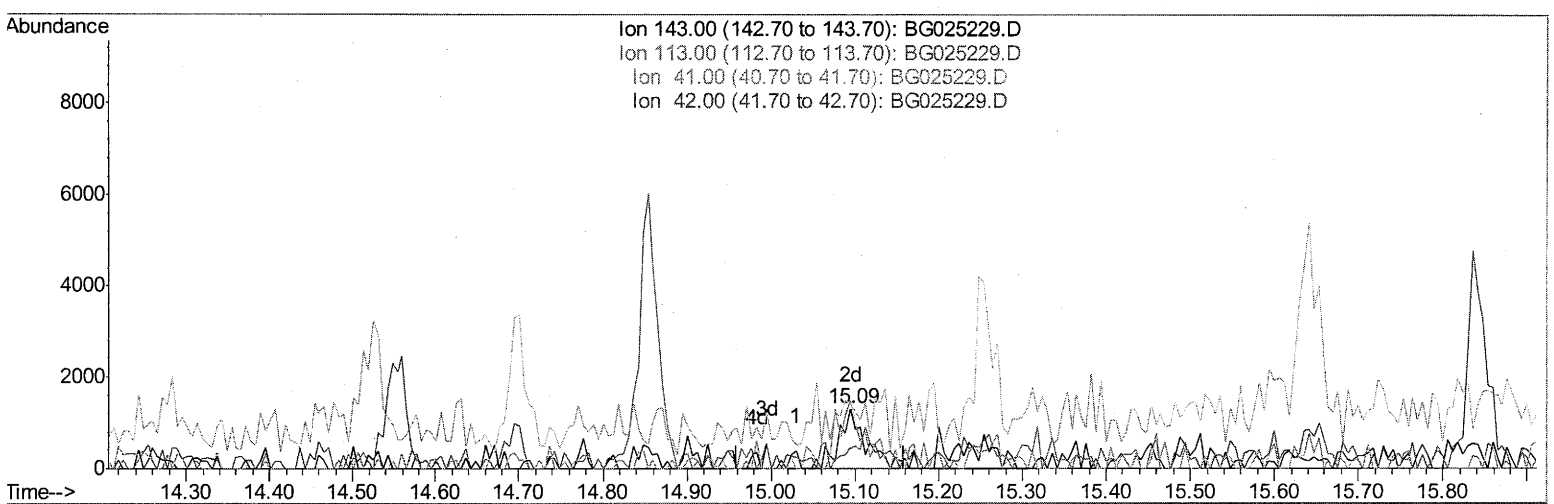
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Instrument :  
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Manual Integrations  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 16 05:54:49 2016  
 Response via : Initial Calibration



(51) 4-Nitrophenol-d4 (S)

15.095min (+0.035) 0.92ng/ul m

response 2704

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 143.00 | 100    | 100     |
| 113.00 | 117.30 | 113.87  |
| 41.00  | 58.00  | 108.32# |
| 42.00  | 42.20  | 34.75   |

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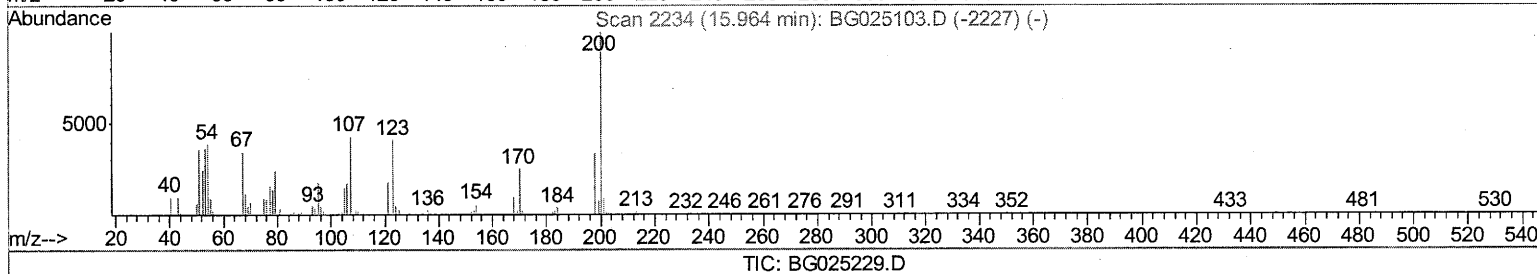
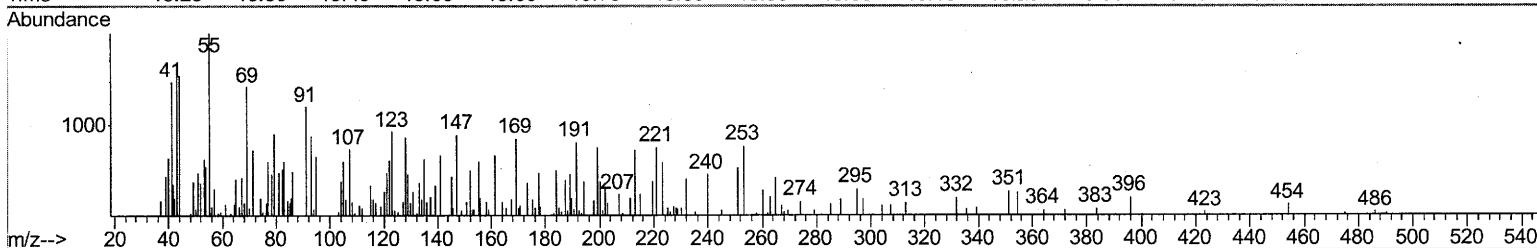
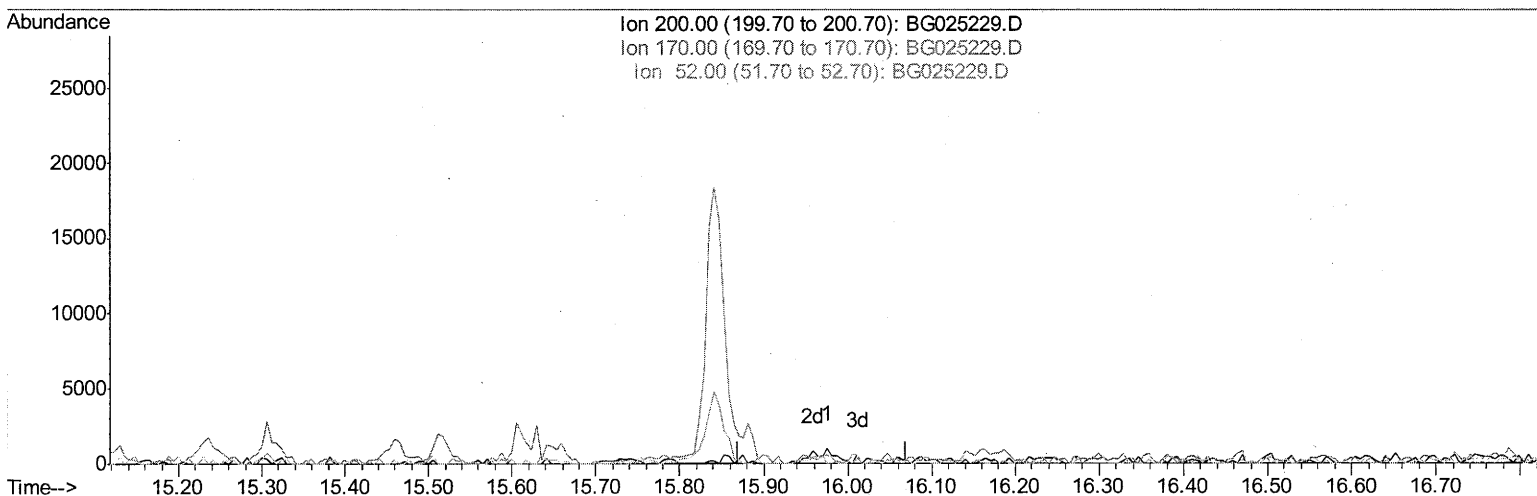
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Data File : BG025229.D  
Acq On : 16 Dec 2016 2:29  
Operator : UM/SJ  
Sample : H5995-16  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA884

Manual Integrations  
APPROVED

UMANGI  
12/16/2016 6:21:22 PM

Quant Time: Dec 16 14:12:56 2016  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Dec 16 05:54:49 2016  
Response via : Initial Calibration



(62) 4,6-Dinitro-2-methylphenol-d2 (S)

15.970min (-15.970) 0.00ng/ul

response 0

| Ion    | Exp%  | Act% |
|--------|-------|------|
| 200.00 | 100   | 0.00 |
| 170.00 | 23.20 | 0.00 |
| 52.00  | 47.20 | 0.00 |
| 0.00   | 0.00  | 0.00 |

# Quantitation Report (Qedit)

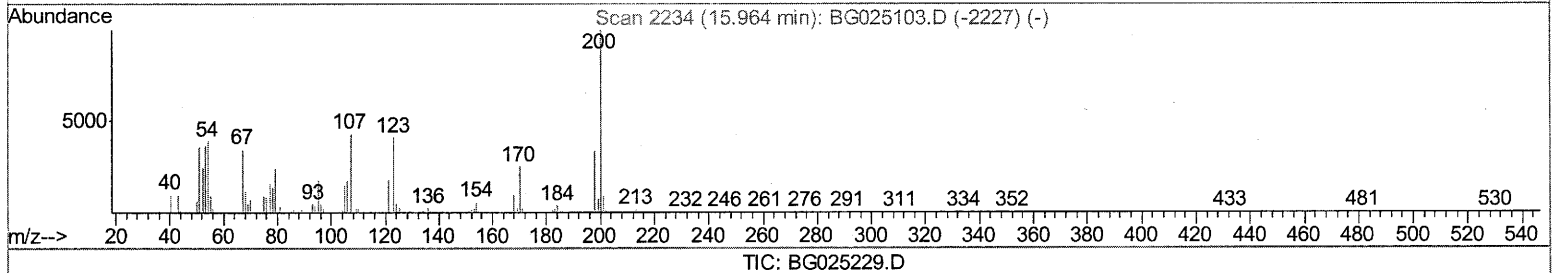
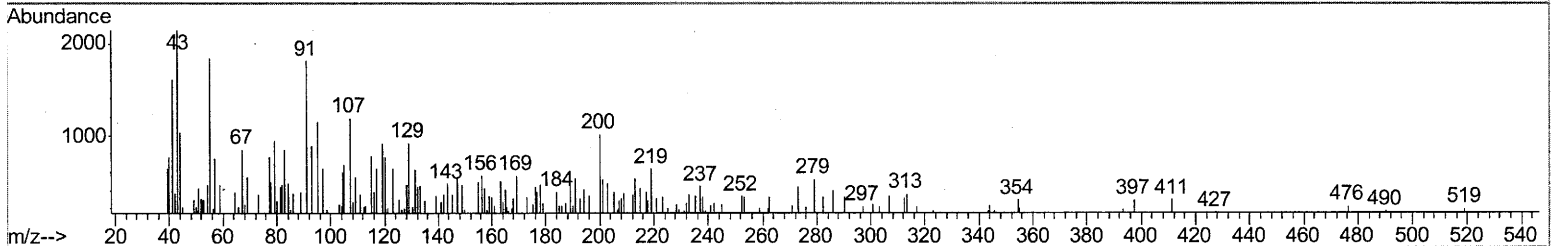
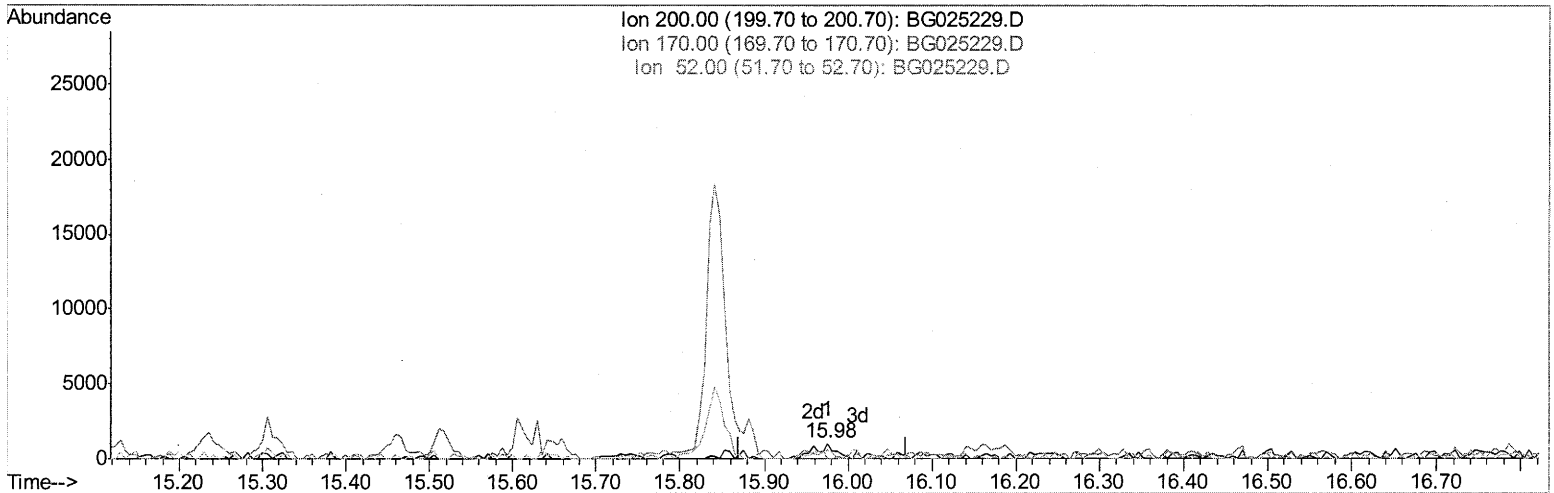
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 Data File : BG025229.D  
 Acq On : 16 Dec 2016 2:29  
 Operator : UM/SJ  
 Sample : H5995-16  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampled :  
 DA884

Manual Integrations  
 APPROVED

UMANGI  
 12/16/2016 6:21:22 PM

Quant Time: Dec 16 14:12:56 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 16 05:54:49 2016  
 Response via : Initial Calibration



(62) 4,6-Dinitro-2-methylphenol-d2 (S)

15.976min (+0.006) 0.60ng/ul m

response 1840

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 200.00 | 100   | 100    |
| 170.00 | 23.20 | 0.00#  |
| 52.00  | 47.20 | 30.97# |
| 0.00   | 0.00  | 0.00   |

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 12/16/16

Data Path : Z:\HPCHEM1\BNA\_G\Data\BG121516\  
 Data File : BG025229.D  
 Acq On : 16 Dec 2016 2:29  
 Operator : UM/SJ  
 Sample : H5995-16  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampled :  
 DA884

Manual Integrations  
 APPROVED

UMANGI  
 12/16/2016 6:21:22 PM

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 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.23  | 152  | 70795    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.06 | 136  | 313410   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.85 | 164  | 249027   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.60 | 188  | 492605   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.89 | 240  | 591258   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.32 | 264  | 608736   | 20.00 | ng/ul | 0.01     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.59  | 96  | 3056   | 2.23  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.39  | 99  | 68084  | 11.15 | ng/ul | 0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.55  | 67  | 53966  | 14.28 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.77  | 132 | 37857  | 8.86  | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.94  | 113 | 61904  | 12.13 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.41  | 128 | 31778  | 14.32 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 12714  | 4.60  | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.68 | 165 | 29839  | 5.61  | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.21 | 131 | 69730  | 13.33 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.25 | 166 | 249674 | 14.58 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.55 | 160 | 312598 | 16.66 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.09 | 143 | 2704m  | 0.92  | ng/ul | 0.04 |
| 57) Fluorene-d10               | 15.84 | 176 | 238118 | 14.77 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.98 | 200 | 1840m  | 0.60  | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.70 | 188 | 339080 | 16.31 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.97 | 212 | 399912 | 16.42 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.08 | 264 | 406981 | 16.50 | ng/ul | 0.00 |

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## Target Compounds

|                                |       |     |        |             | Qvalue |
|--------------------------------|-------|-----|--------|-------------|--------|
| 6) Phenol                      | 7.42  | 94  | 16039  | 2.56 ng/ul# | 87     |
| 34) 2-Methylnaphthalene        | 12.70 | 142 | 16425  | 1.43 ng/ul# | 52     |
| 44) Dimethylphthalate          | 14.30 | 163 | 44544  | 2.65 ng/ul# | 97     |
| 69) Phenanthrene               | 17.64 | 178 | 34563  | 1.47 ng/ul  | 97     |
| 74) Fluoranthene               | 19.64 | 202 | 59001  | 2.11 ng/ul# | 95     |
| 77) Pyrene                     | 20.00 | 202 | 63095  | 2.22 ng/ul# | 95     |
| 80) Benzo(a)anthracene         | 21.87 | 228 | 51168  | 1.67 ng/ul# | 88     |
| 81) Bis(2-ethylhexyl)phthalate | 21.72 | 149 | 675451 | 44.10 ng/ul | 97     |
| 82) Chrysene                   | 21.95 | 228 | 65718  | 2.29 ng/ul# | 87     |
| 85) Benzo(b)fluoranthene       | 24.22 | 252 | 86663  | 2.87 ng/ul# | 59     |
| 88) Benzo(a)pyrene             | 25.15 | 252 | 49075  | 1.69 ng/ul# | 64     |

(#) = qualifier out of range (m) = manual integration (+) = signals summed