

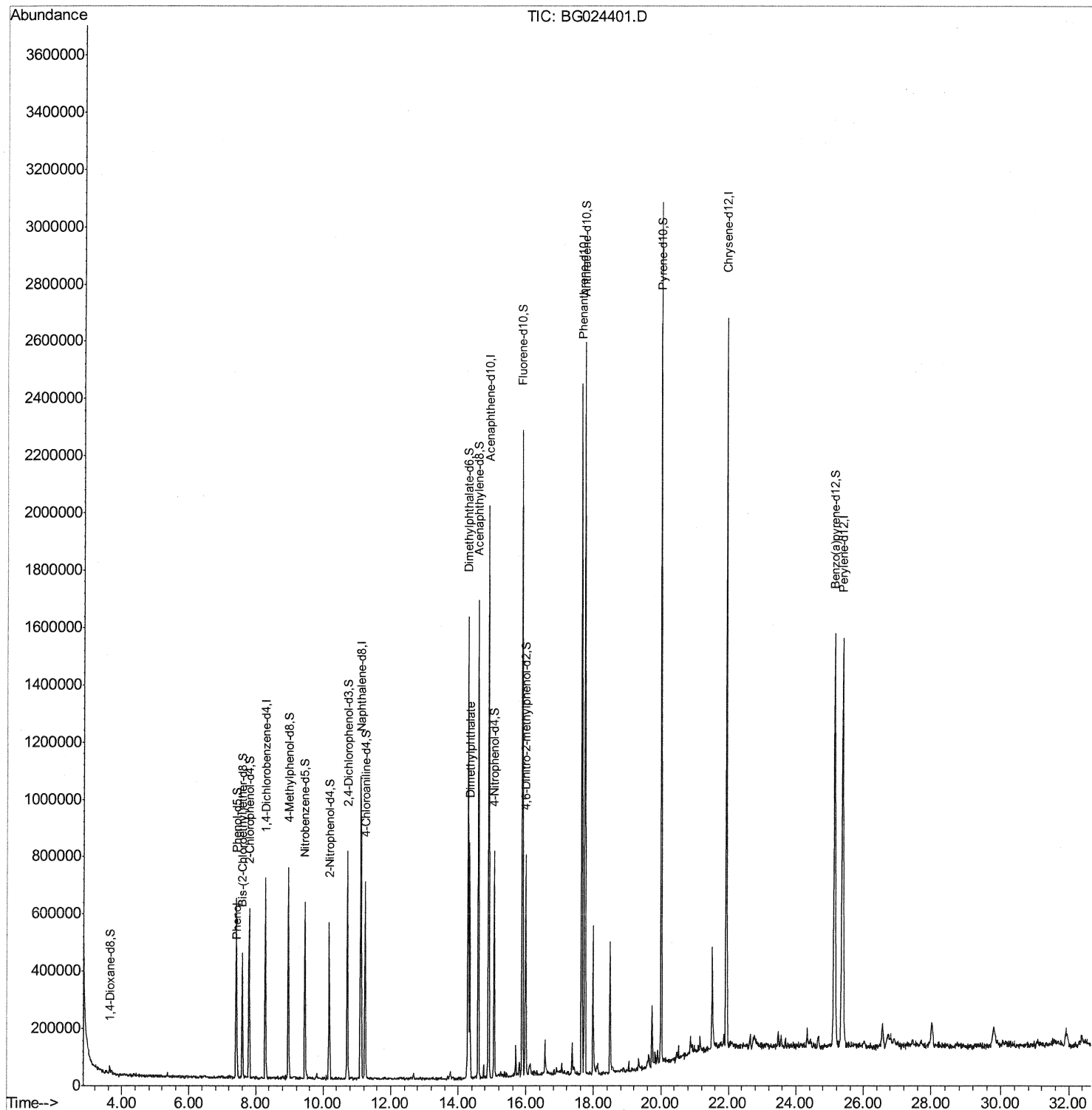
Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG101716\  
 Data File : BG024401.D  
 Acq On : 18 Oct 2016 2:41  
 Operator : UM/SJ  
 Sample : H5241-05  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BDNC5

Manual Integrations  
 APPROVED

umangi  
 10/18/2016 6:11:20 PM

Quant Time: Oct 18 03:45:39 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG0100716.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 18 03:43:57 2016  
 Response via : Initial Calibration



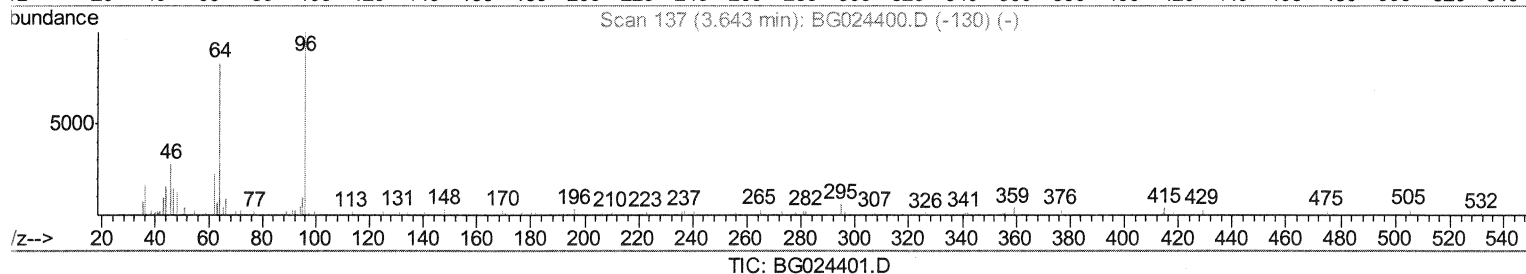
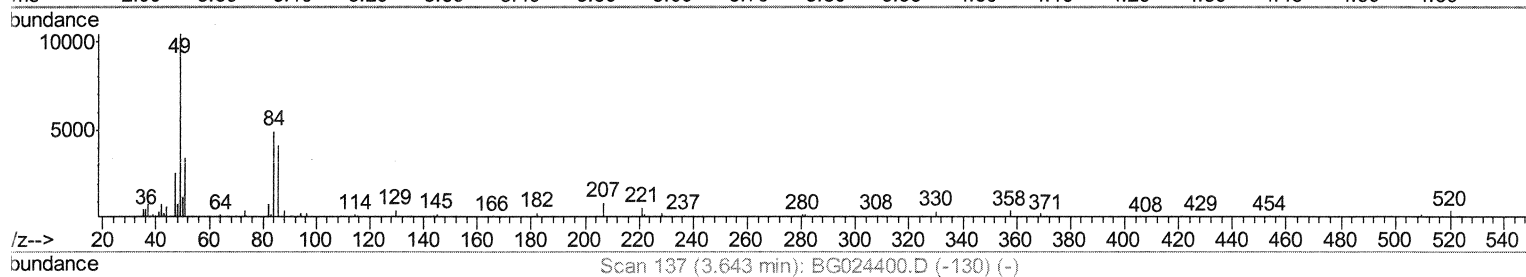
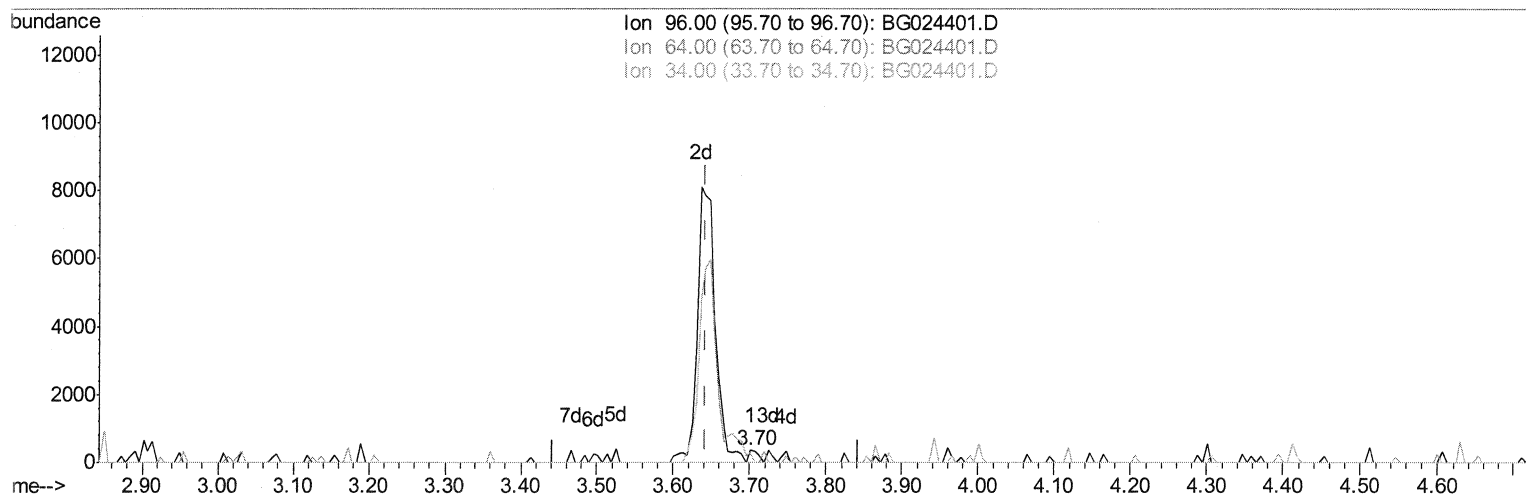
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TIC: BG024401.D

(3) 1,4-Dioxane-d8 (S)

3.701min (+0.058) 0.08ng/uL

response 327

| Ion   | Exp%  | Act%  |
|-------|-------|-------|
| 96.00 | 100   | 100   |
| 64.00 | 88.90 | 77.41 |
| 34.00 | 0.00  | 0.00  |
| 0.00  | 0.00  | 0.00  |

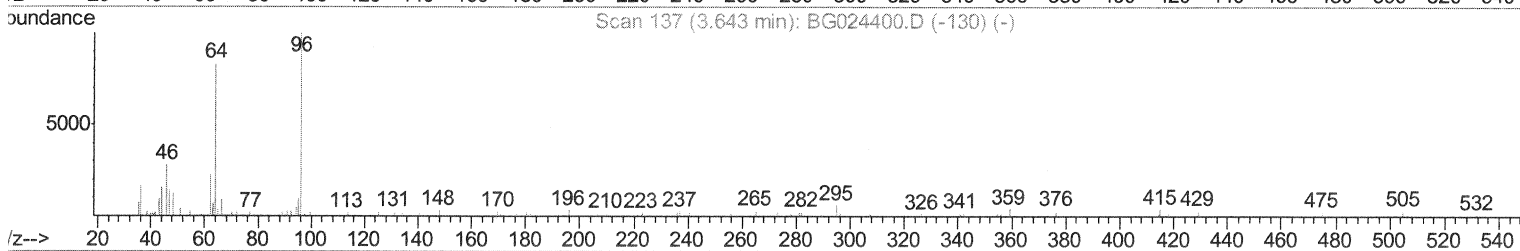
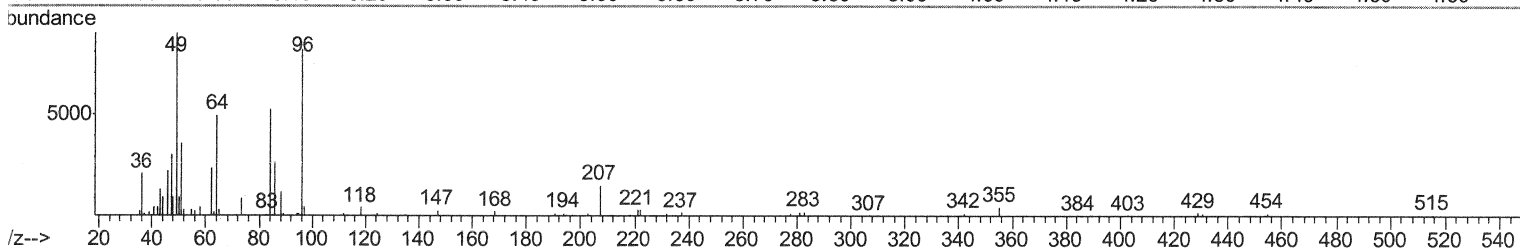
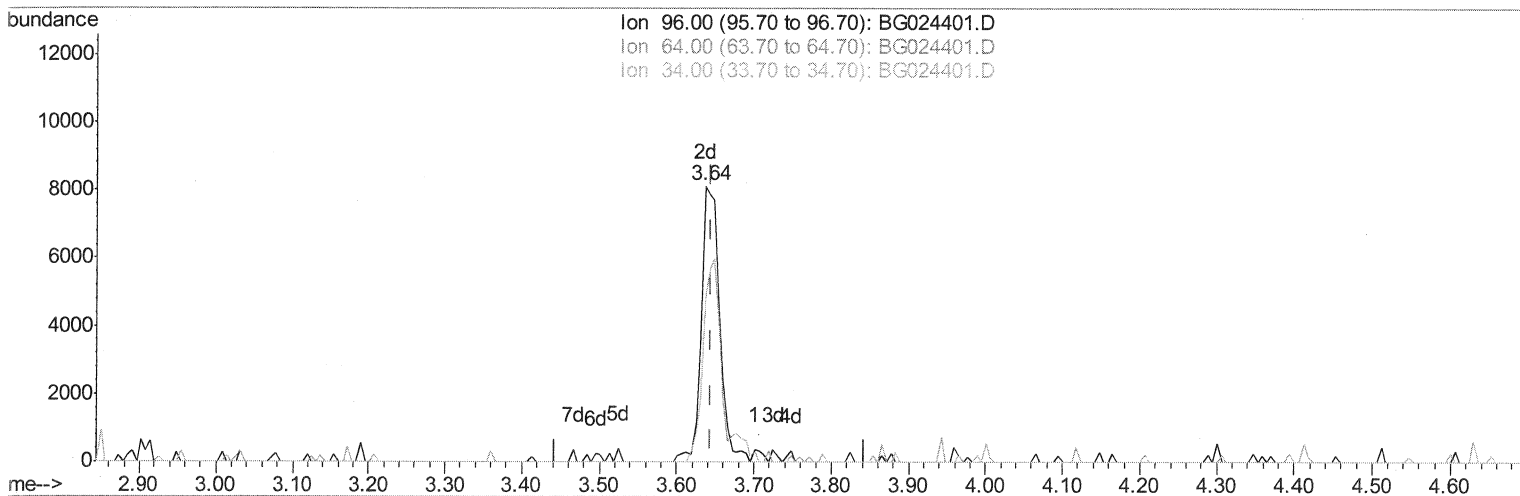
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Manual Integrations  
APPROVED

umangi  
10/18/2016 6:11:20 PM

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Response via : Initial Calibration



TIC: BG024401.D

(3) 1,4-Dioxane-d8 (S)

3.636min (-0.007) 3.49ng/uL m

SJ 10/20/16

response 13457

| Ion   | Exp%  | Act%   |
|-------|-------|--------|
| 96.00 | 100   | 100    |
| 64.00 | 88.90 | 61.07# |
| 34.00 | 0.00  | 0.00   |
| 0.00  | 0.00  | 0.00   |

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG101716\  
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 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
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Manual Integrations  
 APPROVED

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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.28  | 152  | 183291   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.11 | 136  | 845403   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.90 | 164  | 718983   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.64 | 188  | 1491766  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.94 | 240  | 1681076  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.38 | 264  | 1697534  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.64  | 96  | 13457m  | 3.49  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.41  | 99  | 334188  | 19.54 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.60  | 67  | 227109  | 19.65 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.81  | 132 | 239265  | 20.32 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.97  | 113 | 277497  | 20.07 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.45  | 128 | 126842  | 21.14 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.18 | 143 | 154904  | 23.08 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.72 | 165 | 290286  | 19.86 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.24 | 131 | 330750  | 21.53 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.29 | 166 | 1045568 | 20.83 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.59 | 160 | 1204227 | 20.94 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.06 | 143 | 162508  | 18.26 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.88 | 176 | 983047  | 20.71 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.99 | 200 | 176665  | 19.71 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.74 | 188 | 1475779 | 21.90 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 20.01 | 212 | 1632141 | 20.80 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.14 | 264 | 1699942 | 22.26 | ng/ul | 0.00 |

SJ 10/20/16

## Target Compounds

|                       |       |     |        |       |       | Qvalue |
|-----------------------|-------|-----|--------|-------|-------|--------|
| 6) Phenol             | 7.43  | 94  | 27504  | 1.58  | ng/ul | 98     |
| 44) Dimethylphthalate | 14.34 | 163 | 501711 | 10.19 | ng/ul | 100    |

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