

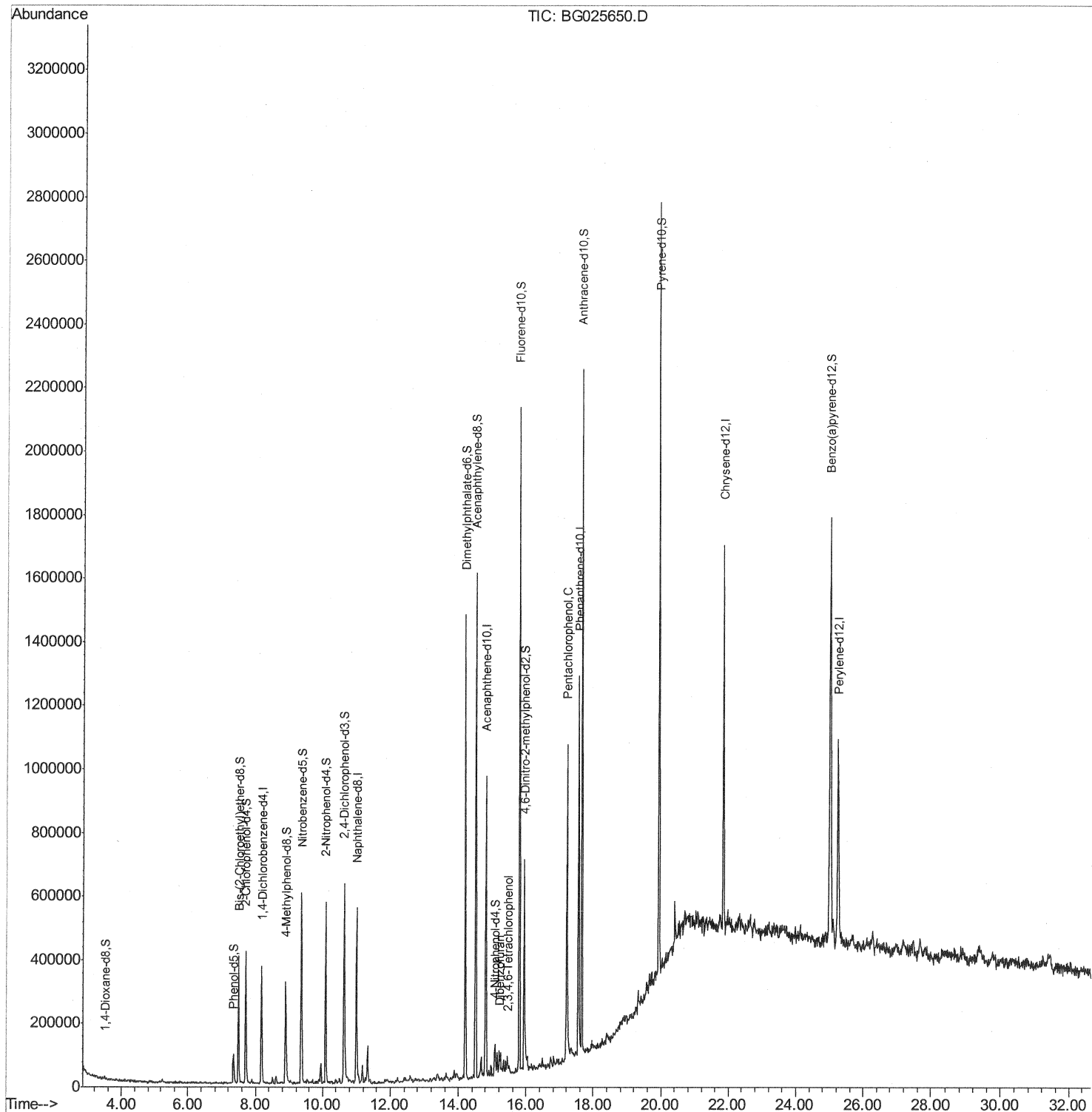
Data Path : Z:\HPCHEM1\BNA\_G\Data\BG012517\  
 Data File : BG025650.D  
 Acq On : 26 Jan 2017 2:14  
 Operator : SJ/MA  
 Sample : I1387-01  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA9Q4

Manual Integrations  
 APPROVED

mohammad  
 1/26/2017 5:11:39 PM

Quant Time: Jan 26 09:58:32 2017  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG011817.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 25 18:08:33 2017  
 Response via : Initial Calibration



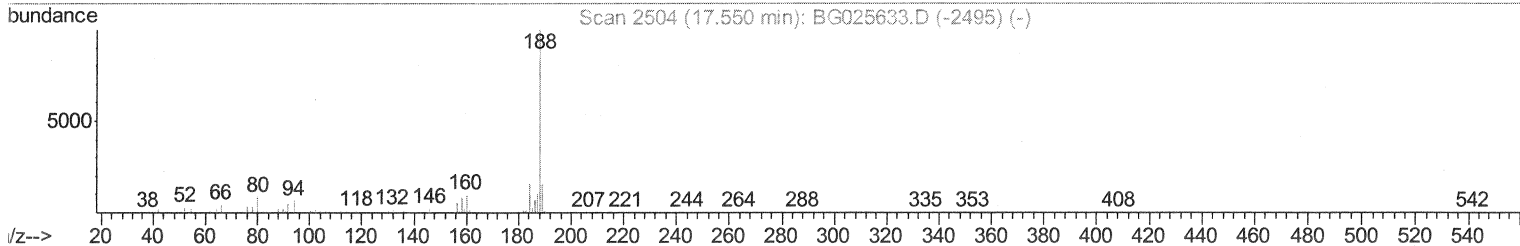
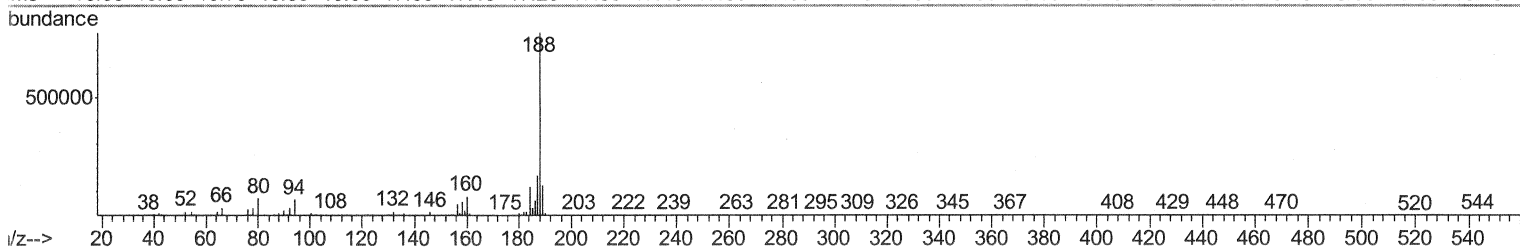
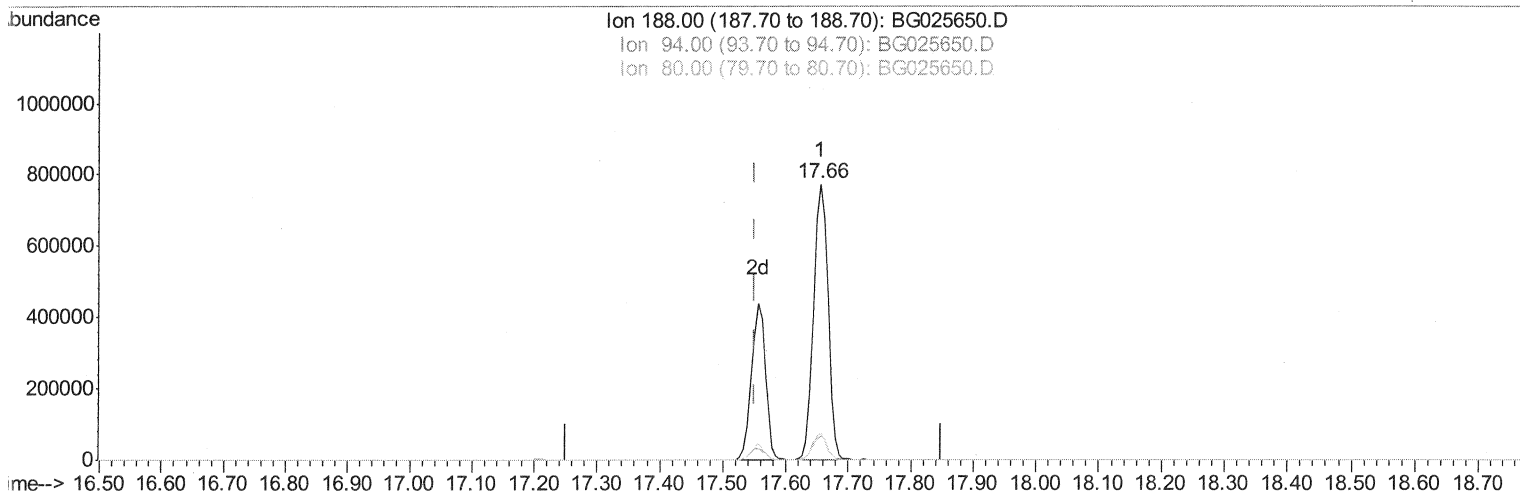
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TIC: BG025650.D

(61) Phenanthrene-d10 (I)

17.655min (+0.105) 20.00ng/ul

response 1235726

| Ion    | Exp% | Act% |
|--------|------|------|
| 188.00 | 100  | 100  |
| 94.00  | 8.80 | 8.81 |
| 80.00  | 9.50 | 9.48 |
| 0.00   | 0.00 | 0.00 |

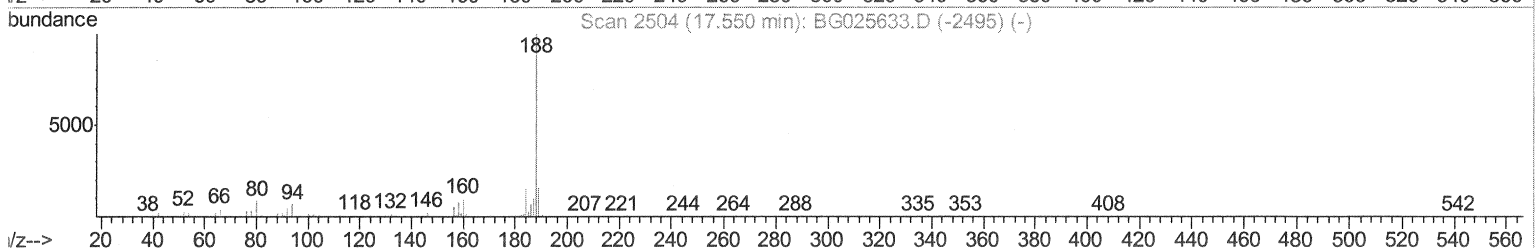
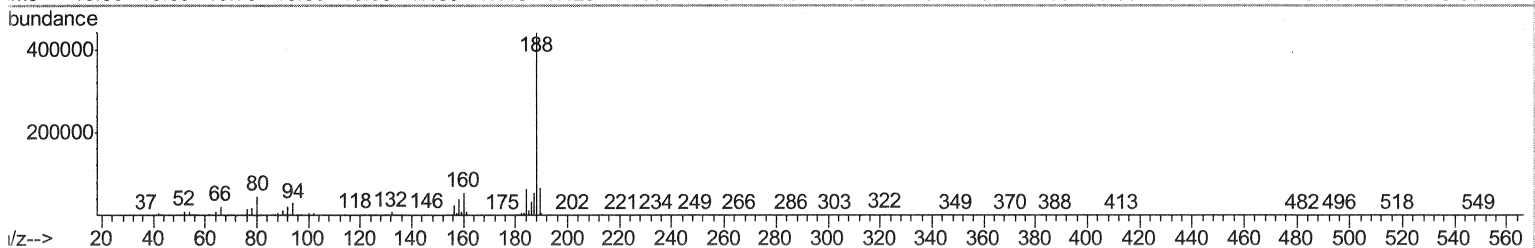
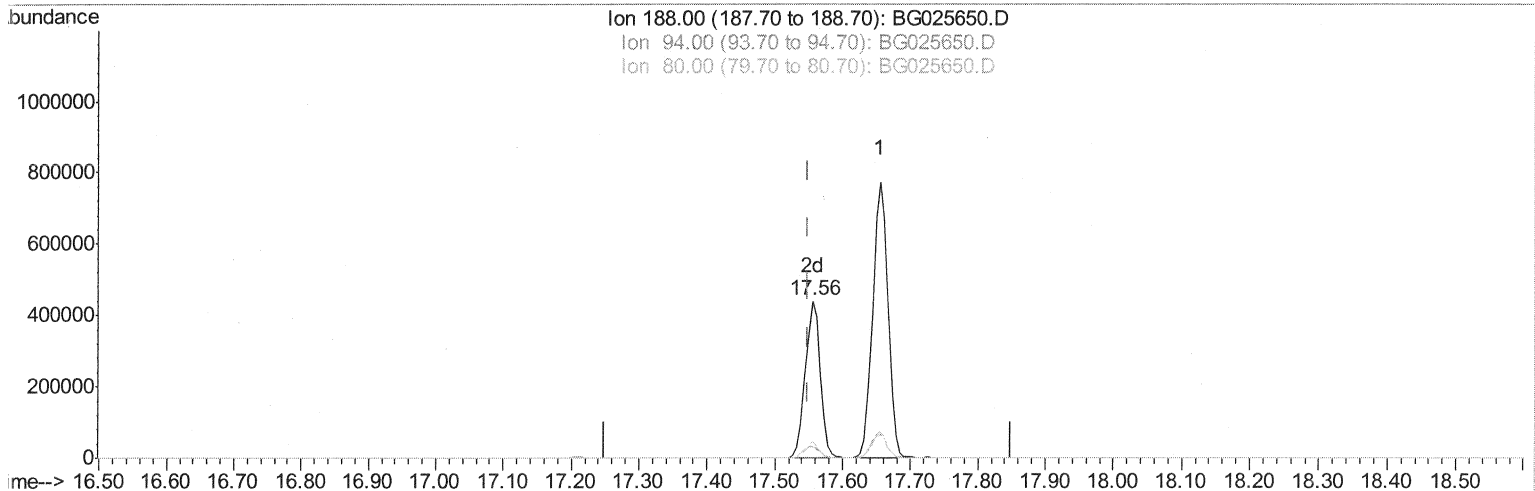
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TIC: BG025650.D

(61) Phenanthrene-d10 (I)

17.555min (+0.005) 20.00ng/ul m

response 679343

| Ion    | Exp% | Act%  |
|--------|------|-------|
| 188.00 | 100  | 100   |
| 94.00  | 8.80 | 6.88# |
| 80.00  | 9.50 | 10.43 |
| 0.00   | 0.00 | 0.00  |

*SJ*  
 01/27/17

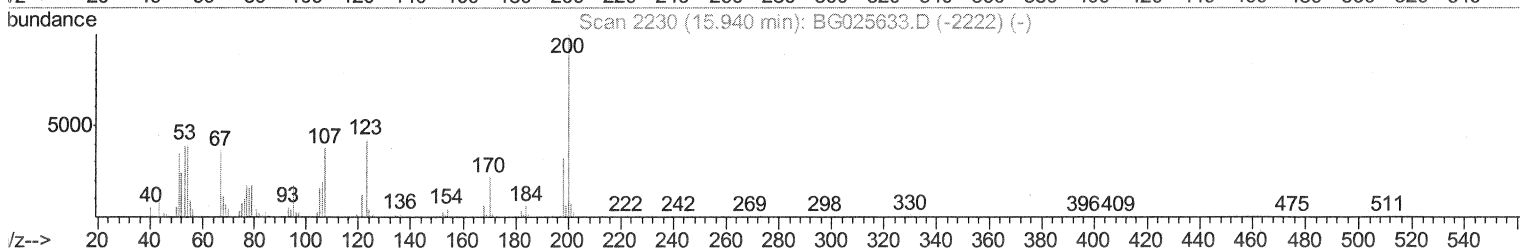
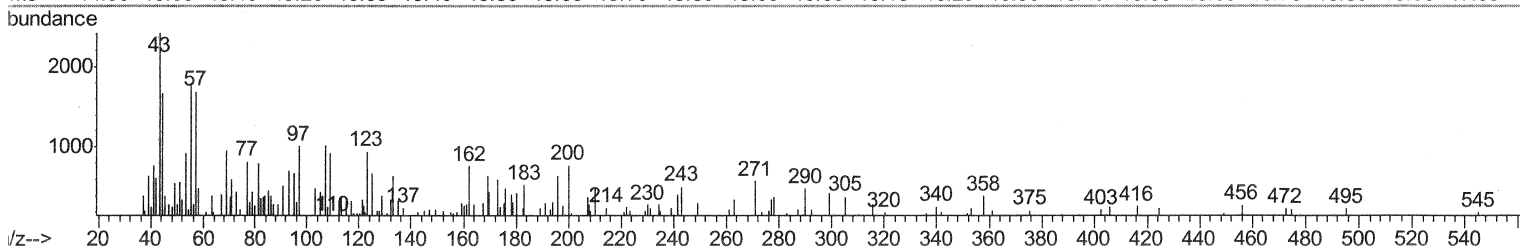
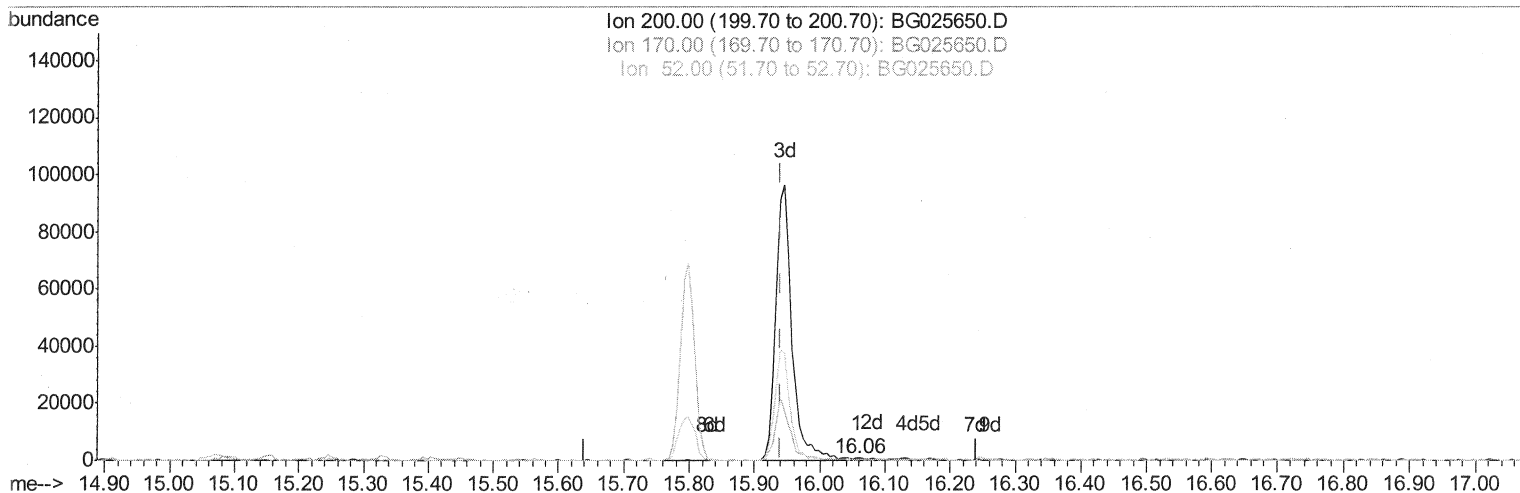
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ALS Vial : 18 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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Manual Integrations  
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mohammad  
1/26/2017 5:11:39 PM

Quant Time: Jan 26 09:57:55 2017  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG011817.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Jan 25 18:08:33 2017  
Response via : Initial Calibration



TIC: BG025650.D

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

16.057min (+0.117) 0.09ng/ul

response 357

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 200.00 | 100   | 100   |
| 170.00 | 23.20 | 0.00# |
| 52.00  | 47.20 | 45.13 |
| 0.00   | 0.00  | 0.00  |

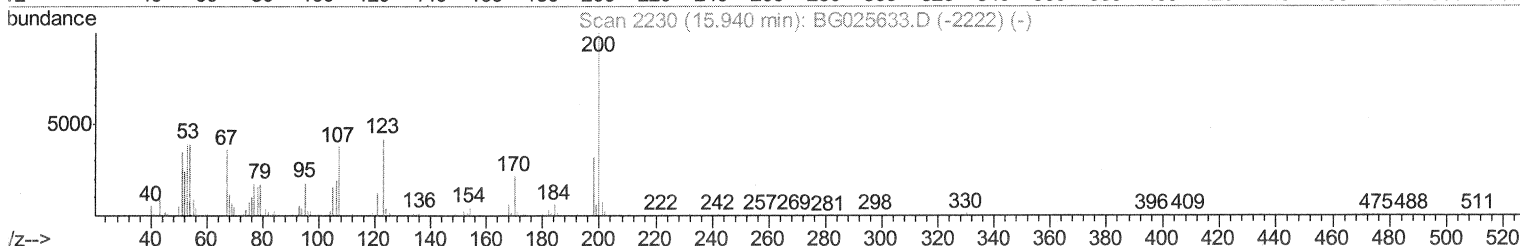
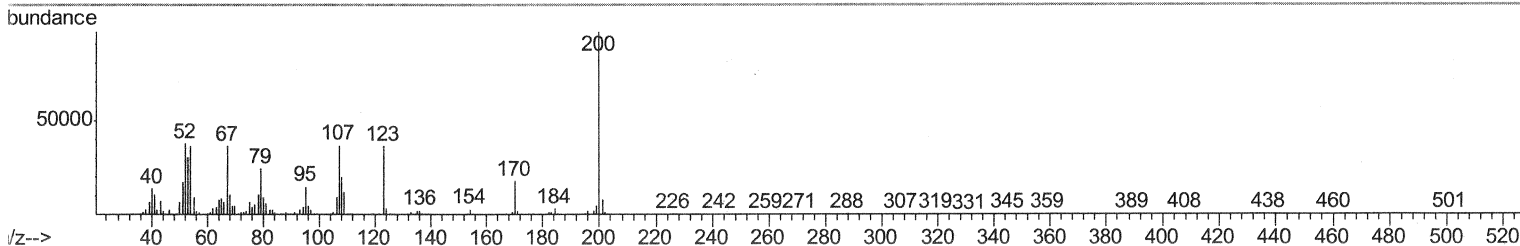
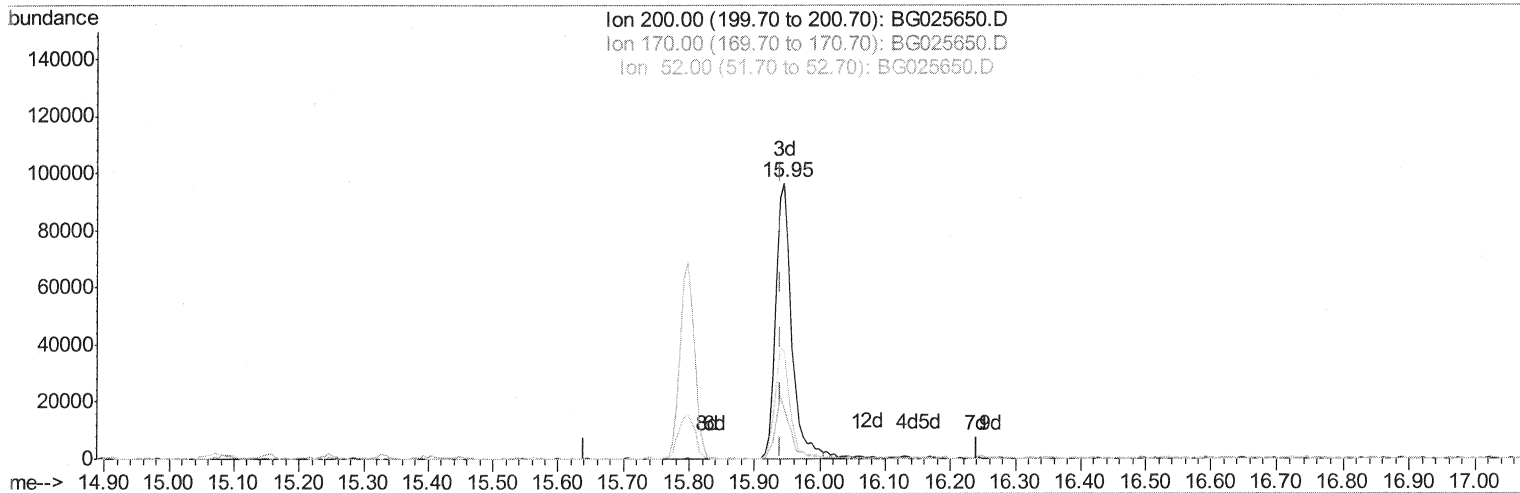
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Operator : SJ/MA  
Sample : I1387-01  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA9Q4

Manual Integrations  
APPROVED

mohammad  
1/26/2017 5:11:39 PM

Quant Time: Jan 26 09:57:55 2017  
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TIC: BG025650.D

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

15.945min (+0.005) 39.26ng/ul m

response 164518

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 200.00 | 100   | 100    |
| 170.00 | 23.20 | 18.37# |
| 52.00  | 47.20 | 38.88  |
| 0.00   | 0.00  | 0.00   |

SJ  
01/27/17

Data Path : Z:\HPCHEM1\BNA\_G\Data\BG012517\  
 Data File : BG025650.D  
 Acq On : 26 Jan 2017 2:14  
 Operator : SJ/MA  
 Sample : I1387-01  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA9Q4

Manual Integrations  
 APPROVED

mohammad  
 1/26/2017 5:11:39 PM

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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.18  | 152  | 91717    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.00 | 136  | 424244   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.81 | 164  | 344043   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.56 | 188  | 679343m  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.85 | 240  | 763360   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.23 | 264  | 793187   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.51  | 96  | 2937    | 1.60  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.34  | 99  | 63799   | 8.12  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.49  | 67  | 213304  | 42.02 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.71  | 132 | 172168  | 31.05 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.89  | 113 | 136801  | 20.47 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.36  | 128 | 121730  | 40.48 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.09 | 143 | 150547  | 41.96 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.63 | 165 | 260740  | 36.83 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.18 | 131 | 4553    | 0.58  | ng/ul | 0.03 |
| 43) Dimethylphthalate-d6       | 14.20 | 166 | 936056  | 38.53 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.51 | 160 | 1091568 | 41.59 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.08 | 143 | 17925   | 5.38  | ng/ul | 0.03 |
| 57) Fluorene-d10               | 15.80 | 176 | 848999  | 37.07 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.95 | 200 | 164518m | 39.26 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.66 | 188 | 1237423 | 42.78 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.93 | 212 | 1357732 | 42.74 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.01 | 264 | 1429927 | 42.89 | ng/ul | 0.02 |

| Target Compounds              |       |     |        |       | Qvalue    |
|-------------------------------|-------|-----|--------|-------|-----------|
| 53) Dibenzofuran              | 15.21 | 168 | 40494  | 1.45  | ng/ul 96  |
| 55) 2,3,4,6-Tetrachlorophenol | 15.45 | 232 | 12703  | 1.83  | ng/ul# 67 |
| 68) Pentachlorophenol         | 17.21 | 266 | 209410 | 56.19 | ng/ul 91  |

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